

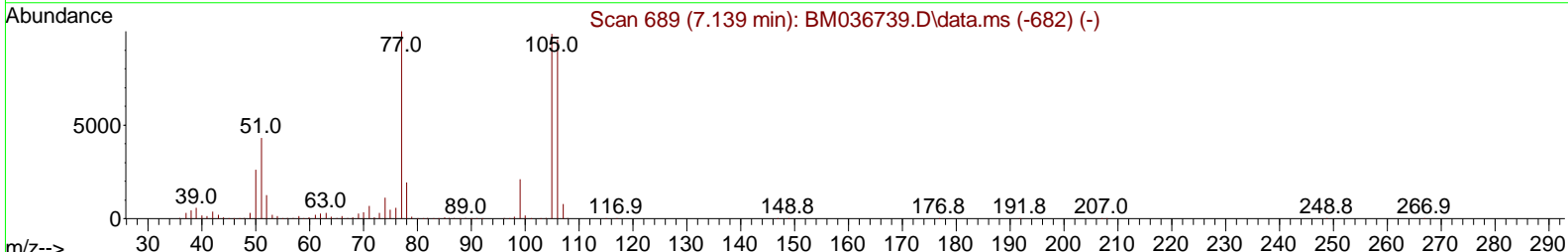
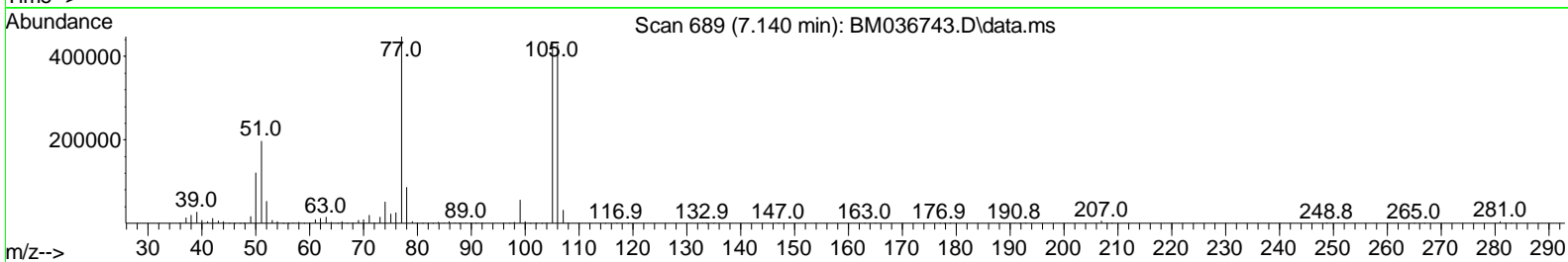
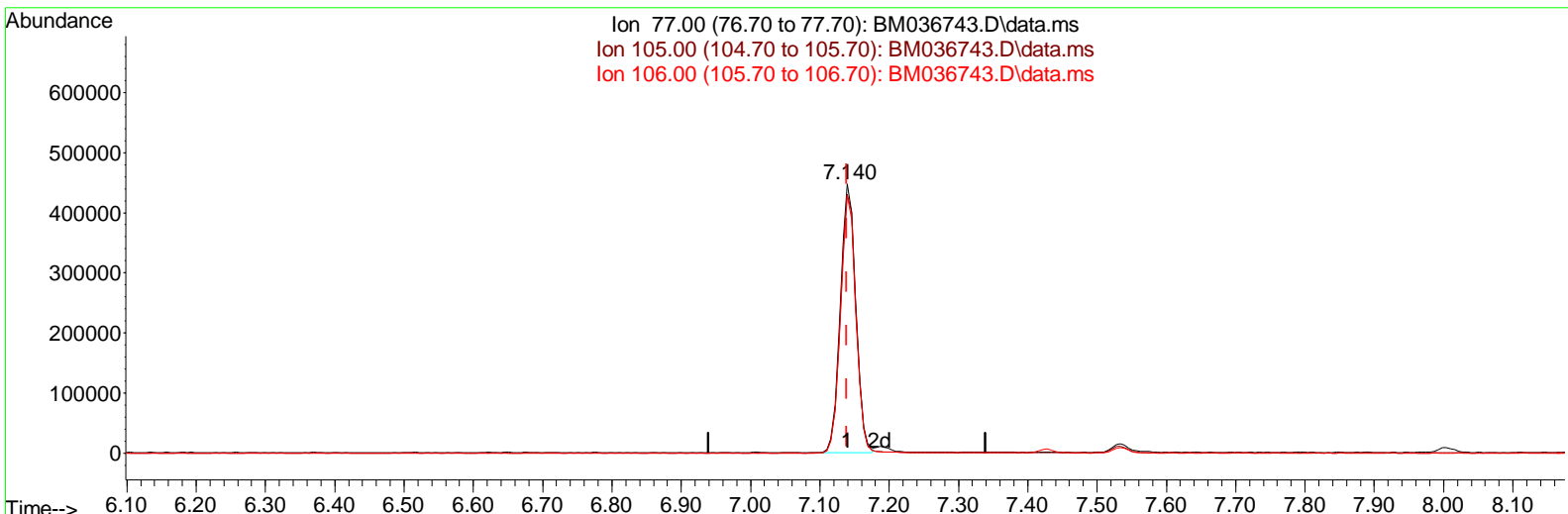
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092622\
 Data File : BM036743.D
 Acq On : 26 Sep 2022 13:42
 Operator : CG/JU
 Sample : PB147814BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS814

Manual Integrations APPROVED

Reviewed By : Jagrut Upadhyay 09/27/2022
 Supervised By : mohammad ahmed 09/28/2022

Quant Time: Sep 26 23:12:50 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Sep 26 23:10:22 2022
 Response via : Initial Calibration



TIC: BM036743.D\data.ms

(6) Benzaldehyde

7.140min (+ 0.000) 42.34 ng/ul

response 686580

Ion	Exp%	Act%
77.00	100.00	100.00
105.00	93.90	96.20
106.00	91.30	95.75
0.00	0.00	0.00

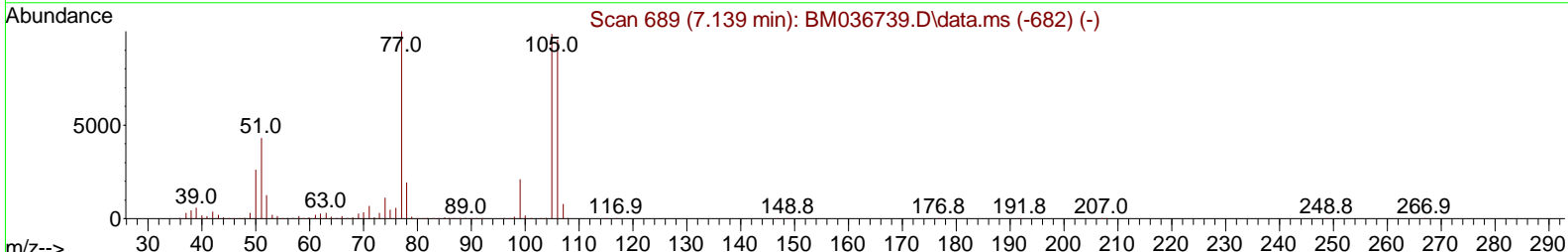
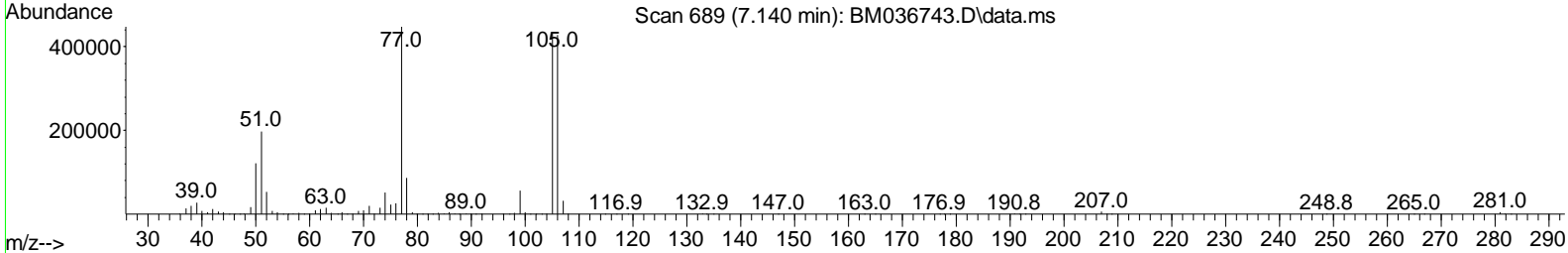
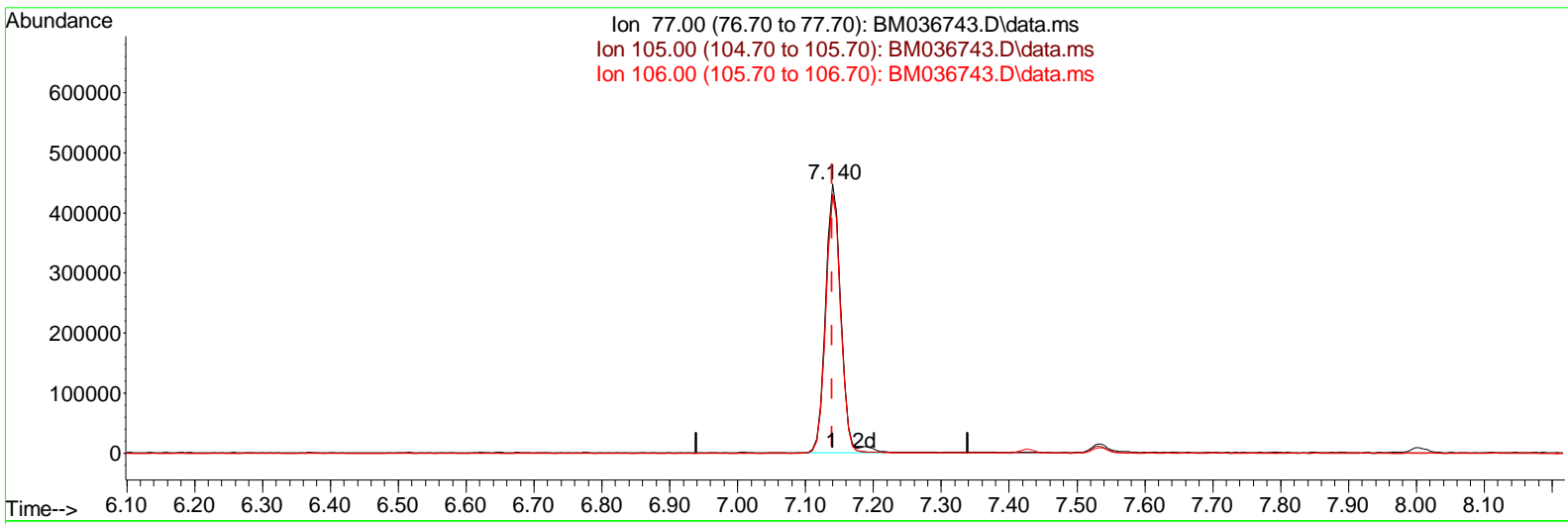
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TIC: BM036743.D\data.ms

(6) Benzaldehyde

7.140min (+ 0.000) 43.33 ng/ul m

response 702590

Ion	Exp%	Act%
77.00	100.00	100.00
105.00	93.90	96.20
106.00	91.30	95.75
0.00	0.00	0.00

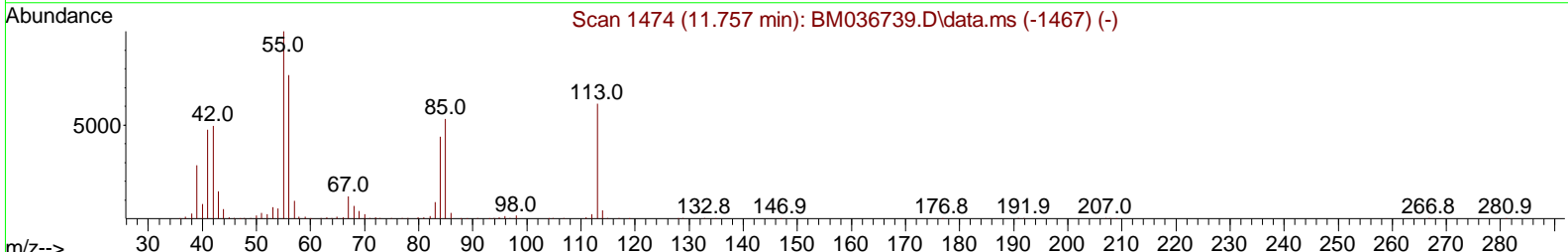
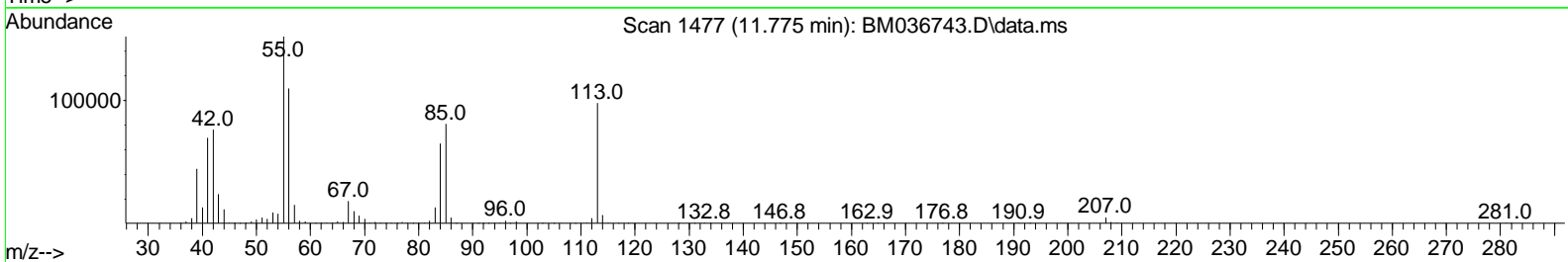
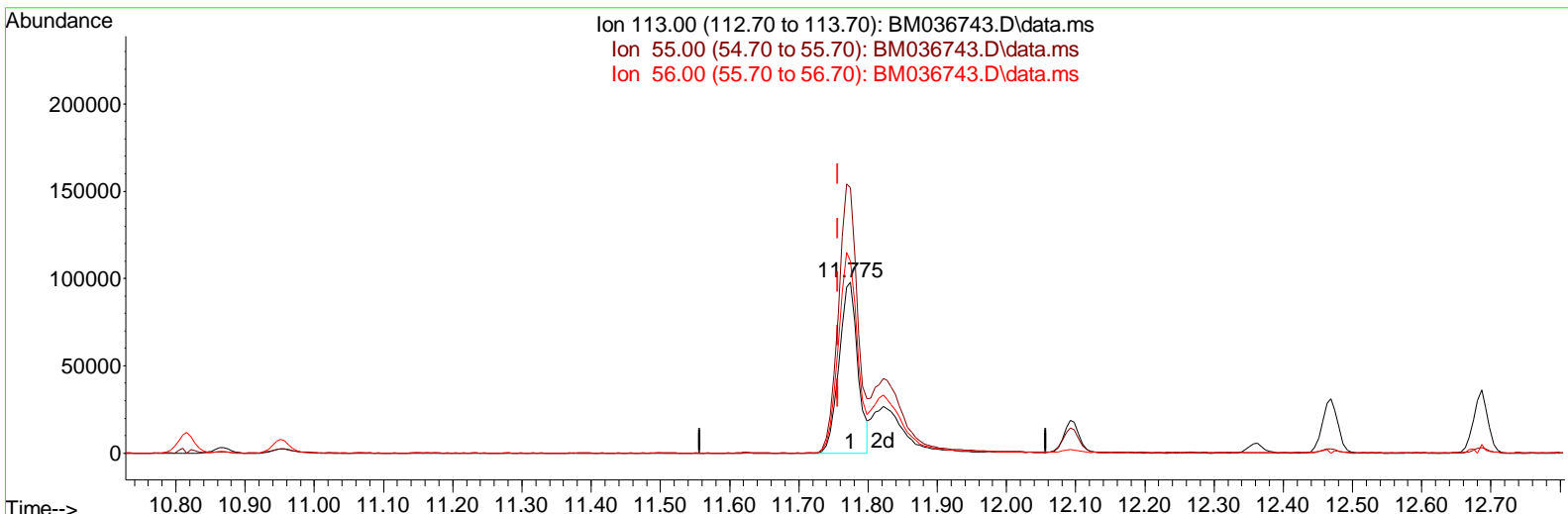
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092622\
 Data File : BM036743.D
 Acq On : 26 Sep 2022 13:42
 Operator : CG/JU
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TIC: BM036743.D\data.ms

(34) Caprolactam

11.775min (+ 0.018) 23.09 ng/ul

response 182342

Ion	Exp%	Act%
113.00	100.00	100.00
55.00	159.40	155.45
56.00	118.30	112.46
0.00	0.00	0.00

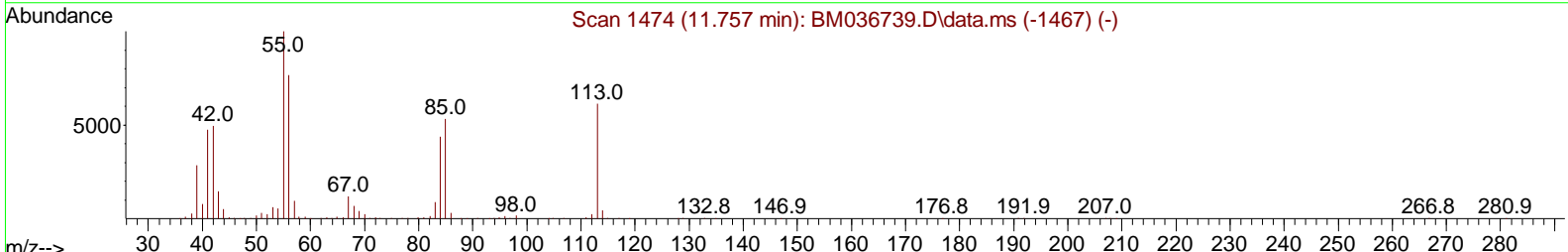
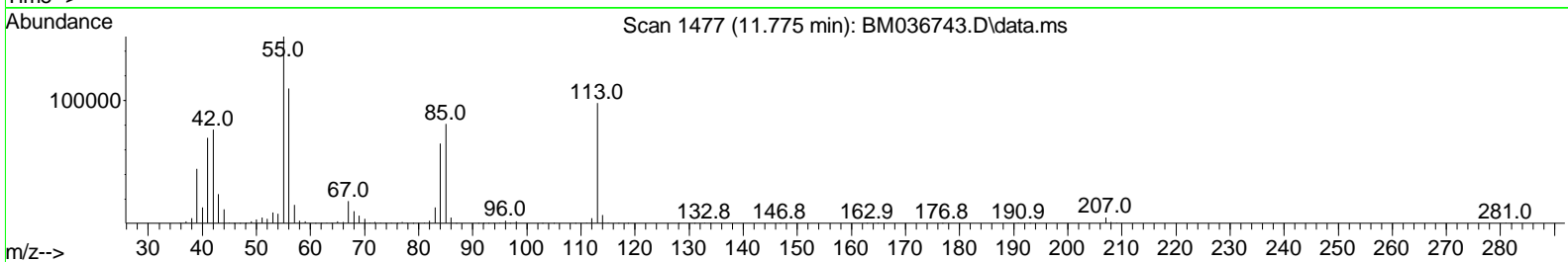
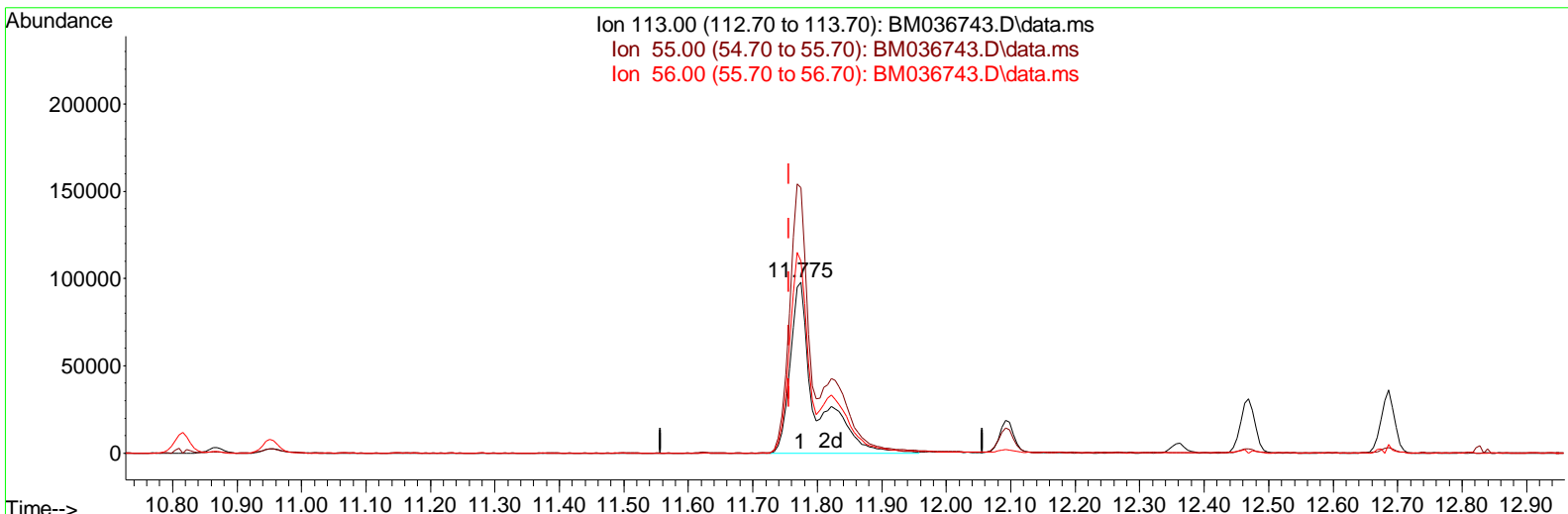
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 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Sep 26 23:10:22 2022
 Response via : Initial Calibration



TIC: BM036743.D\data.ms

(34) Caprolactam

11.775min (+ 0.018) 33.95 ng/ul m

response 268089

Ion	Exp%	Act%
113.00	100.00	100.00
55.00	159.40	155.45
56.00	118.30	112.46
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BMO92622\
 Data File : BMO36743.D
 Acq On : 26 Sep 2022 13:42
 Operator : CG/JU
 Sample : PB147814BS
 Mi sc :
 ALS Vial : 6 Sample Multi plier: 1

Instrument :
 BNA_M
ClientSampleId :
 SLCS814

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 09/27/2022
 Supervised By :mohammad ahmed 09/28/2022

Quant Time: Sep 26 23:12:50 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BMO91522.M
 Quant Title : SVOA CALI BRATI ON
 QLast Update : Mon Sep 26 23:10:22 2022
 Response via : Ini tial Cal i brati on

Compound	R. T.	QI on	Response	Conc	Units	Dev(Mi n)
Internal Standards						
1) 1,4-Di chl orobenzene-d4	8.004	152	374728	20.000	ng/ul	0.00
20) Naphthal ene-d8	10.816	136	1647859	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.627	164	1026251	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.368	188	2039954	20.000	ng/ul	0.00
79) Chrysene-d12	21.533	240	1628573	20.000	ng/ul	0.00
88) Peryl ene-d12	23.968	264	1346382	20.000	ng/ul	0.00
System Moni toring Compounds						
3) 1,4-Di oxane-d8	3.387	96	59676	5.970	ng/uL	0.00
4) Pyri di ne-d5	3.810	84	803076	27.547	ng/ul	0.00
7) Phenol -d5	7.163	99	1181838	35.292	ng/ul	0.00
9) Bi s-(2-Chl oroethyl)eth. . .	7.334	67	718376	35.324	ng/ul	0.00
11) 2-Chl orophenol -d4	7.534	132	904534	37.518	ng/ul	0.00
15) 4-Methyl phenol -d8	8.716	113	911139	36.823	ng/ul	0.00
21) Ni trobenzene-d5	9.175	128	449234	40.277	ng/ul	0.00
24) 2-Ni trophenol -d4	9.892	143	434485	43.490	ng/ul	0.00
28) 2,4-Di chl orophenol -d3	10.434	165	862621	37.258	ng/ul	0.00
31) 4-Chl oroani li ne-d4	10.951	131	1196663	31.087	ng/ul	0.00
46) Di methyl phthal ate-d6	14.045	166	2488547	35.954	ng/ul	0.00
49) Acenaphthyl ene-d8	14.327	160	3030589	36.913	ng/ul	0.00
54) 4-Ni trophenol -d4	14.827	143	456006	33.967	ng/ul	0.00
60) Fl uorene-d10	15.616	176	2223338	35.845	ng/ul	0.00
65) 4,6-Di ni tro-2-methyl ph. . .	15.733	200	328762	40.734	ng/ul	0.00
73) Anthracene-d10	17.468	188	3321607	35.612	ng/ul	0.00
81) Pyrene-d10	19.751	212	3554094	44.360	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.815	264	2471985	37.580	ng/ul	0.01
Target Compounds						
2) 1,4-Di oxane	3.422	88	124145	11.529	ng/uL	98
5) Pyri di ne	3.834	79	883104	30.151	ng/ul	96
6) Benzal dehyde	7.140	77	702590m	43.327	ng/ul	
8) Phenol	7.193	94	1221136	33.976	ng/ul	99
10) Bi s(2-Chl oroethyl)ether	7.428	93	931400	33.307	ng/ul	99
12) 2-Chl orophenol	7.563	128	925315	35.480	ng/ul	98
13) 2-Methyl phenol	8.445	108	887776	33.605	ng/ul	100
14) 2,2'-oxybi s(1-Chl oropr. . .	8.534	45	1124690	33.934	ng/ul	100
16) Acetophenone	8.834	105	1427499	33.260	ng/ul	100
17) N-Ni troso-di -n-propyl a. . .	8.822	70	722929	34.908	ng/ul	97
18) 4-Methyl phenol	8.781	108	985743	34.506	ng/ul	99
19) Hexachl oroethane	9.092	117	361376	34.210	ng/ul	95
22) Ni trobenzene	9.216	77	1078693	36.426	ng/ul	99
23) I sophorone	9.751	82	1970031	32.627	ng/ul	99
25) 2-Ni trophenol	9.928	139	477117	38.666	ng/ul	98
26) 2,4-Di methyl phenol	9.987	107	1001603	32.910	ng/ul	98
27) Bi s(2-Chl oroethoxy)met. . .	10.228	93	1222193	33.960	ng/ul	99
29) 2,4-Di chl orophenol	10.463	162	856182	35.175	ng/ul	99
30) Naphthal ene	10.869	128	2998262	33.400	ng/ul	100
32) 4-Chl oroani li ne	10.975	127	1207818	30.391	ng/ul	100
33) Hexachl orobutadi ene	11.151	225	472507	31.740	ng/ul	99
34) Caprol actam	11.775	113	268089m	33.955	ng/ul	
35) 4-Chl oro-3-methyl phenol	12.092	107	946861	35.240	ng/ul	97

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BMO92622\
 Data File : BMO36743.D
 Acq On : 26 Sep 2022 13: 42
 Operator : CG/JU
 Sample : PB147814BS
 Mi sc :
 ALS Vial : 6 Sample Multi plier: 1

Instrument :
 BNA_M
ClientSampleId :
 SLCS814

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 09/27/2022
 Supervised By :mohammad ahmed 09/28/2022

Quant Time: Sep 26 23: 12: 50 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BMO91522.M
 Quant Title : SVOA CALI BRATI ON
 QLast Update : Mon Sep 26 23: 10: 22 2022
 Response via : Ini tial Cal i brati on

Compound	R. T.	QI on	Response	Conc	Units	Dev(Mi n)
36) 2-Methyl naphthal ene	12. 469	142	2002805	33. 486	ng/ul	98
37) 1-Methyl naphthal ene	12. 686	142	2044456	34. 002	ng/ul	100
39) 1, 2, 4, 5-Tetrachl oroben. . .	12. 833	216	967277	33. 300	ng/ul	100
40) Hexachl orocycl opentadi ene	12. 810	237	432214	28. 552	ng/ul	99
41) 2, 4, 6-Tri chl orophenol	13. 069	196	626310	35. 215	ng/ul	97
42) 2, 4, 5-Tri chl orophenol	13. 139	196	693197	36. 354	ng/ul	100
43) 1, 1' -Bi phenyl	13. 469	154	2556033	33. 787	ng/ul	99
44) 2-Chl oronaphthal ene	13. 510	162	1982510	33. 396	ng/ul	100
45) 2-Ni troani li ne	13. 716	65	603038	41. 272	ng/ul	97
47) Di methyl phthal ate	14. 092	163	2470773	33. 652	ng/ul	99
48) 2, 6-Di ni trotol uene	14. 210	165	481413	40. 330	ng/ul	99
50) Acenaphthyl ene	14. 357	152	3144474	33. 500	ng/ul	99
51) 3-Ni troani li ne	14. 539	138	538590	37. 620	ng/ul #	95
52) Acenaphthene	14. 692	153	2202322	33. 532	ng/ul	100
53) 2, 4-Di ni trophenol	14. 739	184	185538	35. 030	ng/ul	98
55) 4-Ni trophenol	14. 839	109	370378	32. 274	ng/ul	99
56) Di benzofuran	15. 027	168	2959205	33. 550	ng/ul	98
57) 2, 4-Di ni trotol uene	14. 992	165	697817	40. 211	ng/ul	96
58) 2, 3, 4, 6-Tetrachl orophenol	15. 251	232	558340	34. 980	ng/ul #	97
59) Di ethyl phthal ate	15. 451	149	2533784	34. 474	ng/ul	100
61) Fl uorene	15. 674	166	2578612	34. 326	ng/ul	99
62) 4-Chl orophenyl -phenyl e. . .	15. 669	204	1200838	33. 425	ng/ul	99
63) 4-Ni troani li ne	15. 698	138	524269	37. 691	ng/ul	99
66) 4, 6-Di ni tro-2-methyl ph. . .	15. 751	198	332608	36. 519	ng/ul	99
67) N-Ni trosodi phenyl ami ne	15. 880	169	2099020	34. 936	ng/ul	99
68) 4-Bromophenyl -phenyl ether	16. 557	248	675439	33. 593	ng/ul	97
69) Hexachl orobenzene	16. 674	284	766411	32. 227	ng/ul	99
70) Atrazi ne	16. 833	200	694600	32. 761	ng/ul	98
71) Pentachl orophenol	17. 015	266	437841	30. 995	ng/ul	99
72) Phenanthrene	17. 410	178	3887317	34. 596	ng/ul	99
74) Anthracene	17. 504	178	3921780	34. 338	ng/ul	100
75) 1, 2, 3, 4-Tetrachl oroben. . .	13. 433	216	999788	35. 834	ng/uL	99
76) Pentachl orobenzene	14. 945	250	893250	29. 600	ng/uL	99
77) Carbazol e	17. 768	167	3480021	32. 911	ng/ul	100
78) Di -n-butyl phthal ate	18. 333	149	4405019	36. 754	ng/ul	99
80) Fl uoranthene	19. 415	202	4174982	42. 524	ng/ul	99
82) Pyrene	19. 780	202	4322860	41. 596	ng/ul	99
83) Butyl benzyl phthal ate	20. 674	149	1716898	43. 599	ng/ul	95
84) 3, 3' -Di chl orobenzi di ne	21. 450	252	1143165	35. 924	ng/ul	98
85) Benzo(a)anthracene	21. 521	228	3598381	35. 265	ng/ul	100
86) Bi s(2-ethyl hexyl)phtha. . .	21. 450	149	2690638	46. 774	ng/ul	99
87) Chrysene	21. 574	228	3345709	34. 544	ng/ul	99
89) Di -n-octyl phthal ate	22. 386	149	3885269	47. 866	ng/ul	100
90) Benzo(b)fl uoranthene	23. 227	252	3130904	36. 671	ng/ul	100
91) Benzo(k)fl uoranthene	23. 274	252	3120138	36. 573	ng/ul	98
93) Benzo(a)pyrene	23. 862	252	2675722	35. 649	ng/ul	98
94) I ndeno(1, 2, 3-cd)pyrene	26. 509	276	2858125	32. 608	ng/ul	96
95) Di benzo(a, h)anthracene	26. 527	278	2566328	32. 285	ng/ul	99
96) Benzo(g, h, i)peryl ene	27. 285	276	2508904	33. 018	ng/ul	98

(#) = qual i fier out of range (m) = manual i ntegrati on (+) = signal s summed

Instrument :

BNA_M

ClientSampleId :

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