

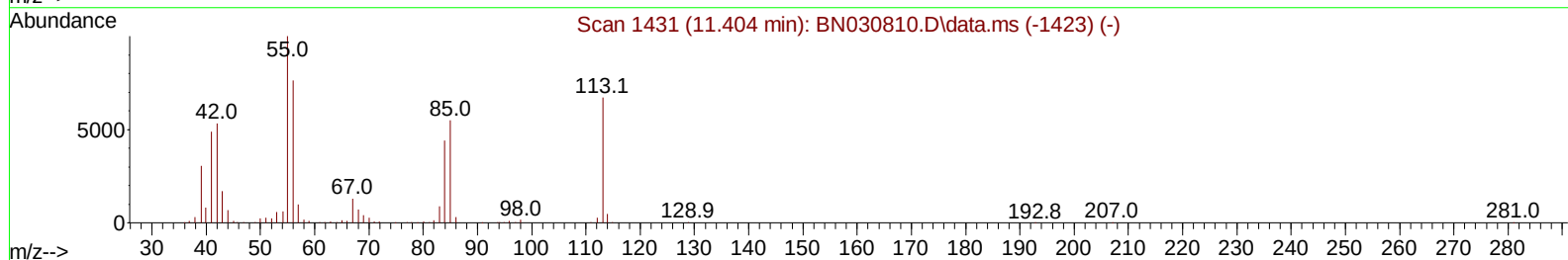
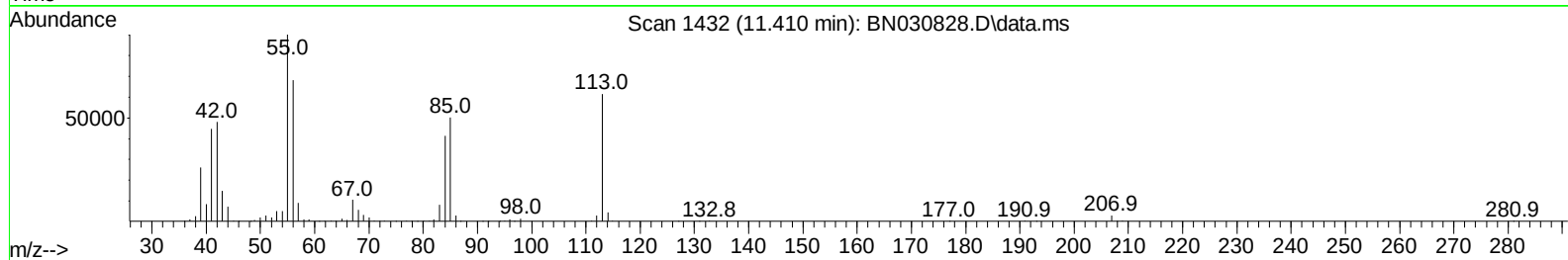
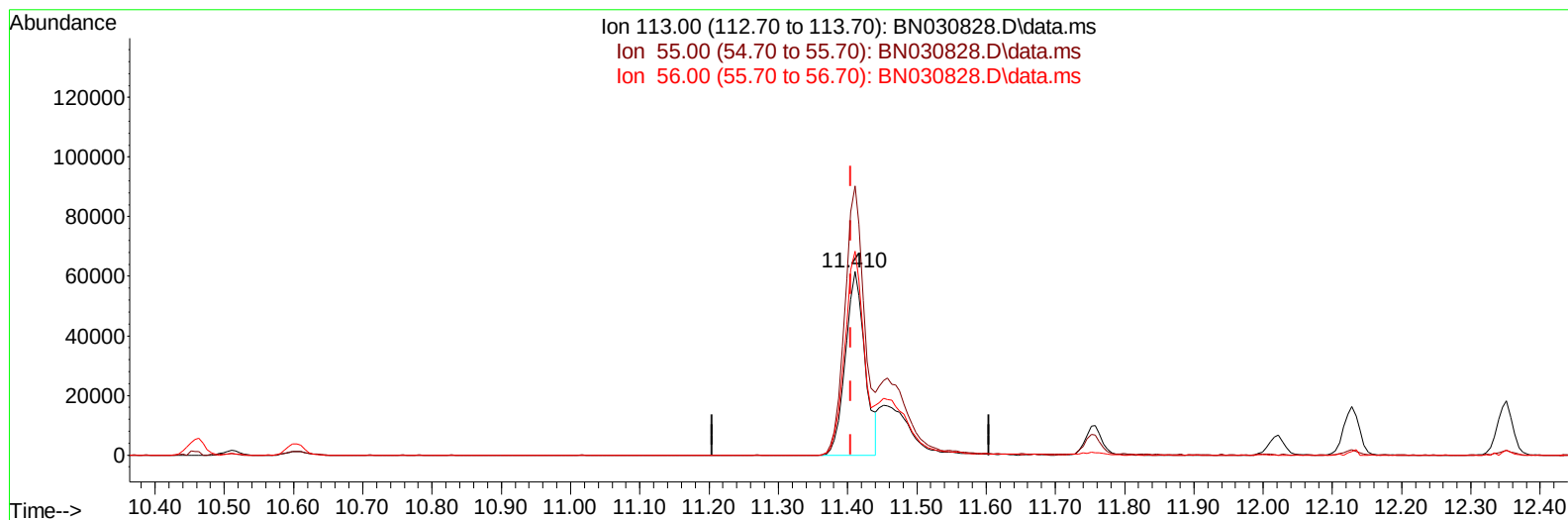
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041024\
Data File : BN030828.D
Acq On : 10 Apr 2024 23:07
Operator : MA/JU
Sample : PB160141BS
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
SLCS141

Manual IntegrationsAPPROVED

Quant Time: Apr 10 23:39:56 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN032824.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Apr 09 12:10:11 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 04/12/2024
Supervised By :mohammad ahmed 04/12/2024



TIC: BN030828.D\data.ms

(34) Caprolactam

11.410min (+ 0.006) 25.67 ng/ul

response 119445

Ion	Exp%	Act%
113.00	100.00	100.00
55.00	155.90	146.63
56.00	113.20	111.14
0.00	0.00	0.00

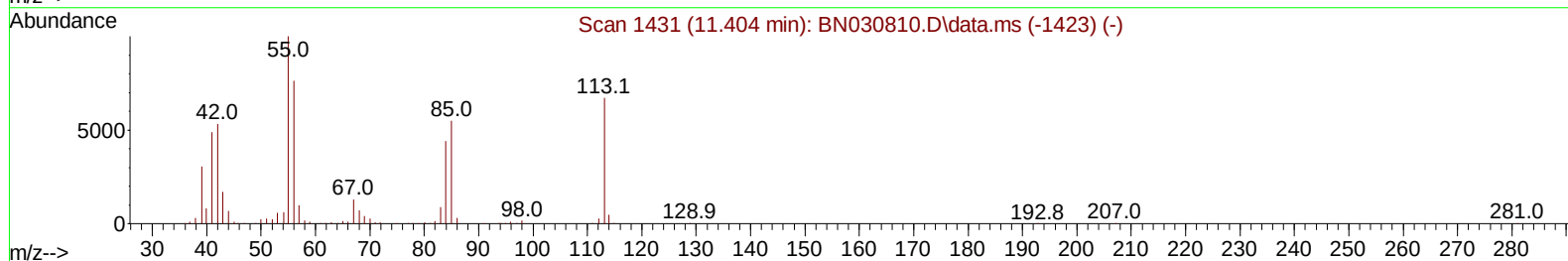
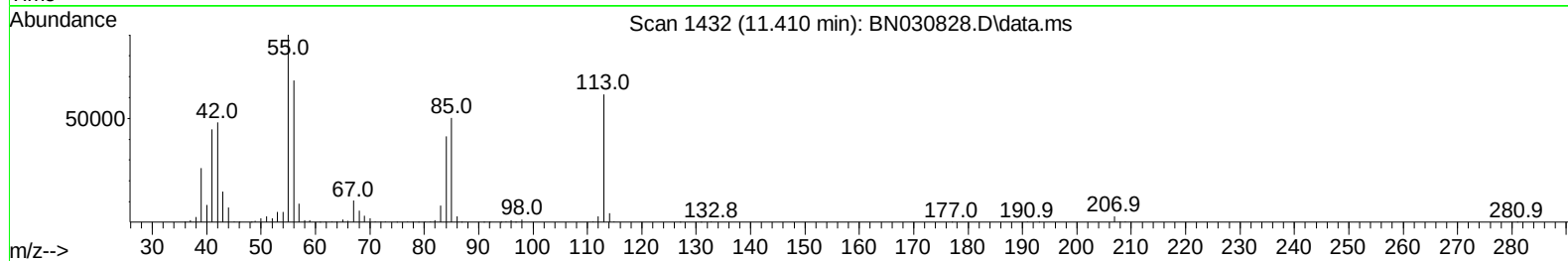
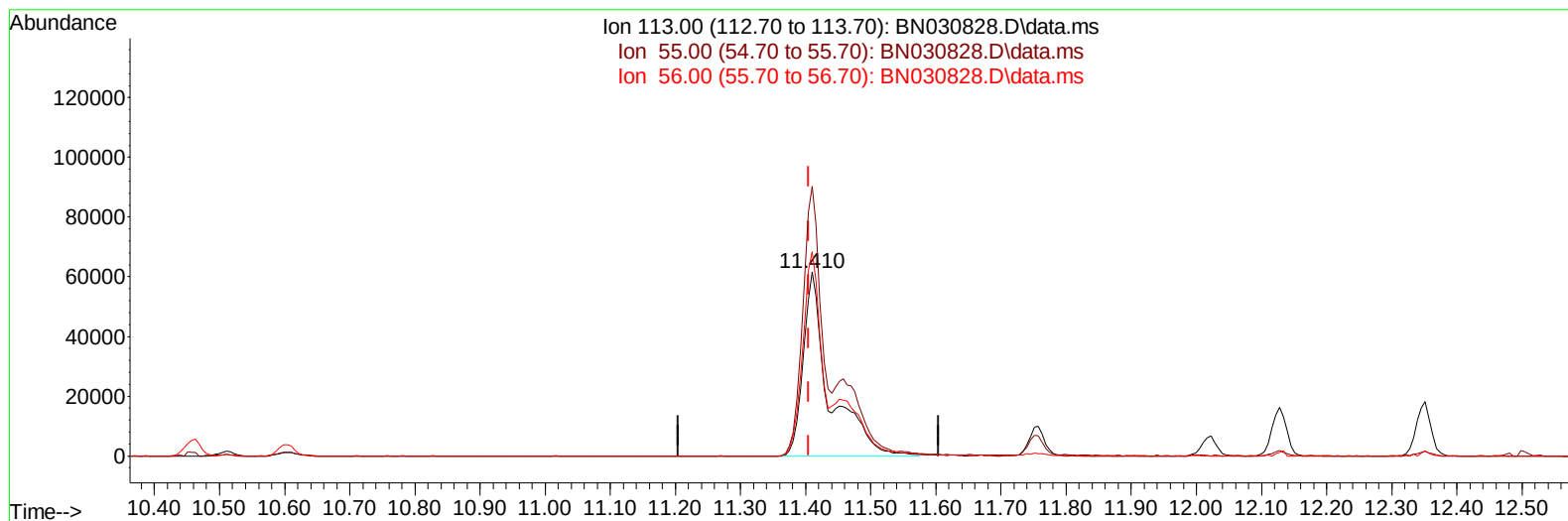
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TIC: BN030828.D\data.ms

(34) Caprolactam

11.410min (+ 0.006) 36.86 ng/ul m

response 171548

Ion	Exp%	Act%
113.00	100.00	100.00
55.00	155.90	146.63
56.00	113.20	111.14
0.00	0.00	0.00

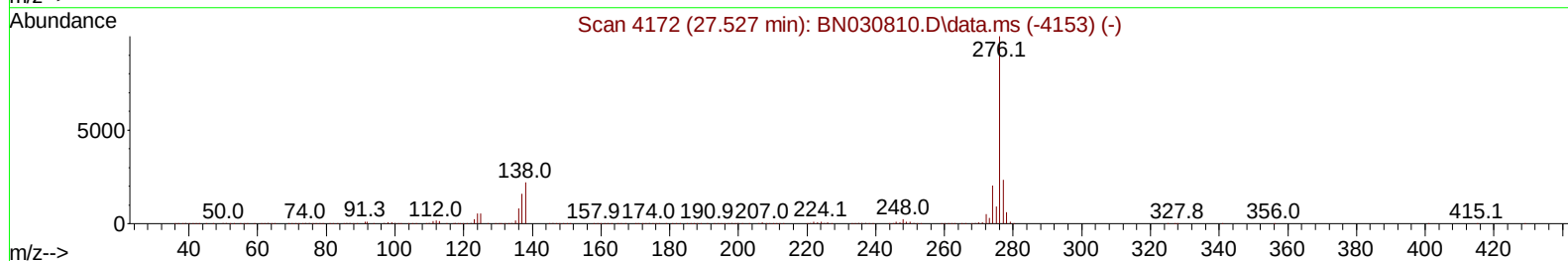
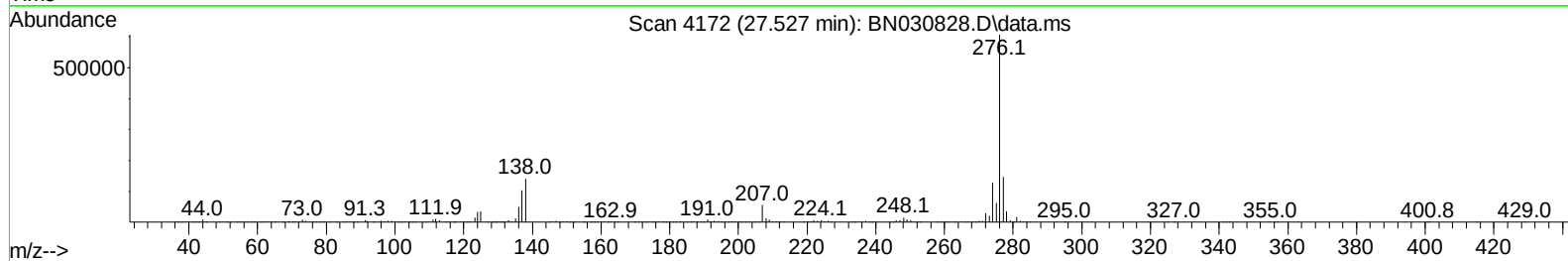
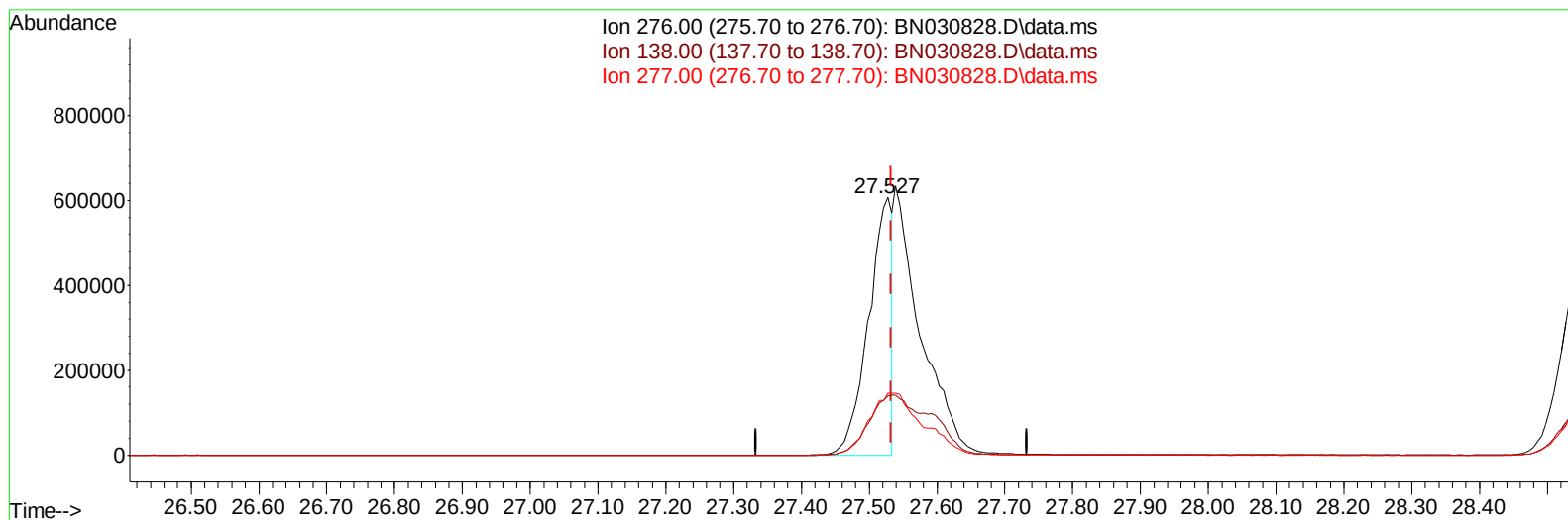
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041024\
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Operator : MA/JU
Sample : PB160141BS
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
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TIC: BN030828.D\data.ms

(94) Indeno(1,2,3-cd)pyrene

27.527min (-0.006) 17.63 ng/u1

response 1472315

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	22.30	23.10
277.00	23.00	24.09
0.00	0.00	0.00

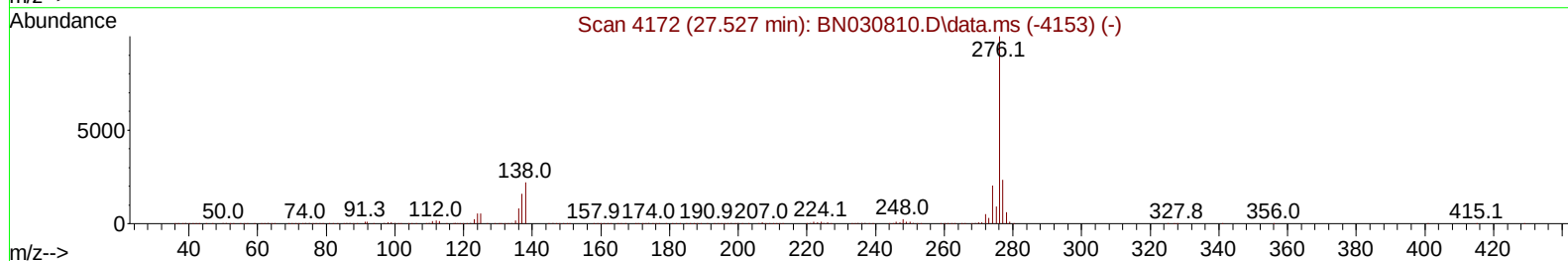
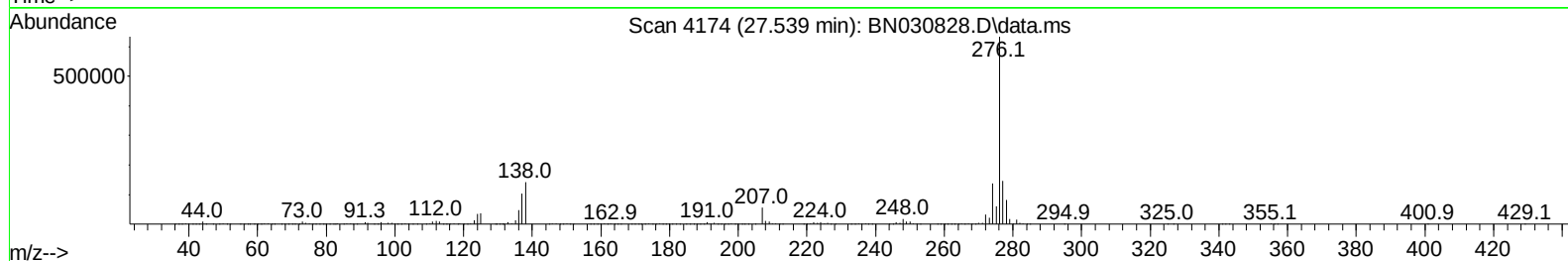
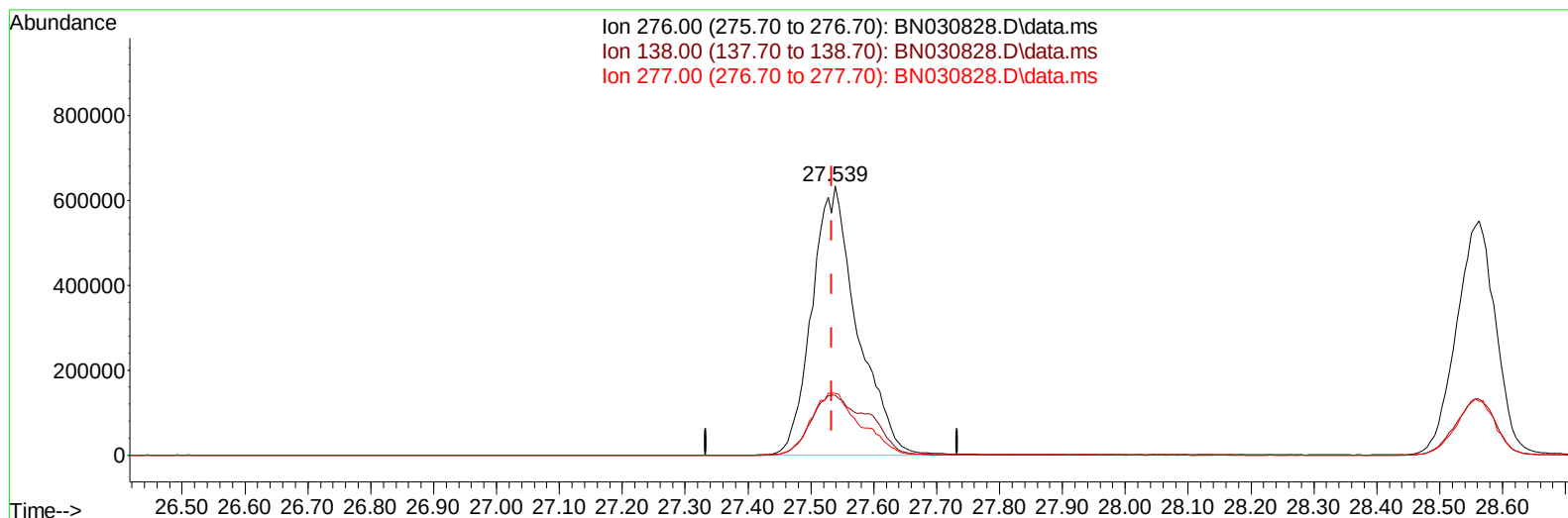
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Quant Title : SVOA CALIBRATION
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TIC: BN030828.D\data.ms

(94) Indeno(1,2,3-cd)pyrene

27.539min (+ 0.006) 38.03 ng/ul m

response 3175101

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	22.30	22.29
277.00	23.00	23.07
0.00	0.00	0.00

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 Sample : PB160141BS
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
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 SLCS141

Manual IntegrationsAPPROVED

Quant Time: Apr 10 23:41:12 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN032824.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Apr 09 12:10:11 2024
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 04/12/2024
 Supervised By :mohammad ahmed 04/12/2024

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.675	152	190811	20.000	ng/ul	0.00
20) Naphthalene-d8	10.458	136	878419	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.322	164	607891	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.069	188	1309596	20.000	ng/ul	0.00
79) Chrysene-d12	21.310	240	1285094	20.000	ng/ul	0.00
88) Perylene-d12	24.239	264	1389289	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.187	96	33907	8.331	ng/uL	0.00
4) Pyridine-d5	3.593	84	423734	34.726	ng/ul	0.00
7) Phenol-d5	6.852	99	582483	37.736	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.017	67	336001	35.243	ng/ul	0.00
11) 2-Chlorophenol-d4	7.211	132	456051	38.947	ng/ul	0.00
15) 4-Methylphenol-d8	8.387	113	477082	36.878	ng/ul	0.00
21) Nitrobenzene-d5	8.834	128	237681	39.006	ng/ul	0.00
24) 2-Nitrophenol-d4	9.552	143	252757	45.342	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.087	165	478991	39.249	ng/ul	0.00
31) 4-Chloroaniline-d4	10.605	131	686426	34.166	ng/ul	0.00
46) Dimethylphthalate-d6	13.740	166	1526769	37.618	ng/ul	0.00
49) Acenaphthylene-d8	14.010	160	1779942	38.143	ng/ul	0.00
54) 4-Nitrophenol-d4	14.534	143	319163	41.112	ng/ul	0.00
60) Fluorene-d10	15.316	176	1320889	37.191	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.445	200	239757	41.477	ng/ul	0.00
73) Anthracene-d10	17.169	188	2123241	37.629	ng/ul	0.00
81) Pyrene-d10	19.469	212	2592308	37.520	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.039	264	2536824	39.629	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.223	88	48149	11.045	ng/uL	98
5) Pyridine	3.611	79	402482	32.688	ng/ul	97
6) Benzaldehyde	6.828	77	292094	40.635	ng/ul	94
8) Phenol	6.875	94	574290	35.714	ng/ul	99
10) Bis(2-Chloroethyl)ether	7.111	93	434312	33.121	ng/ul	99
12) 2-Chlorophenol	7.246	128	457632	36.432	ng/ul	95
13) 2-Methylphenol	8.116	108	431570	34.829	ng/ul	97
14) 2,2'-oxybis(1-Chloropr...	8.199	45	576336	31.214	ng/ul	98
16) Acetophenone	8.499	105	687963	33.682	ng/ul	99
17) N-Nitroso-di-n-propyla...	8.487	70	342660	33.309	ng/ul	96
18) 4-Methylphenol	8.452	108	494495	35.648	ng/ul	96
19) Hexachloroethane	8.746	117	176619	33.537	ng/ul	97
22) Nitrobenzene	8.875	77	515094	34.628	ng/ul	98
23) Isophorone	9.399	82	1001608	33.763	ng/ul	100
25) 2-Nitrophenol	9.581	139	266294	40.647	ng/ul	97
26) 2,4-Dimethylphenol	9.646	107	471348	32.370	ng/ul	99
27) Bis(2-Chloroethoxy)met...	9.881	93	626095	33.195	ng/ul	100
29) 2,4-Dichlorophenol	10.110	162	460216	36.946	ng/ul	94
30) Naphthalene	10.510	128	1527756	33.911	ng/ul	100
32) 4-Chloroaniline	10.628	127	632789	32.681	ng/ul	98
33) Hexachlorobutadiene	10.787	225	264094	33.683	ng/ul	96
34) Caprolactam	11.410	113	171548m	36.864	ng/ul	
35) 4-Chloro-3-methylphenol	11.757	107	526437	37.522	ng/ul	99

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 04/12/2024
 Supervised By :mohammad ahmed 04/12/2024

Quant Time: Apr 10 23:41:12 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN032824.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Apr 09 12:10:11 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
36) 2-Methylnaphthalene	12.128	142	1102712	35.048	ng/ul	98
37) 1-Methylnaphthalene	12.352	142	1087888	34.061	ng/ul	95
39) 1,2,4,5-Tetrachloroben...	12.499	216	557711	35.550	ng/ul	98
40) Hexachlorocyclopentadiene	12.475	237	248414	27.883	ng/ul	99
41) 2,4,6-Trichlorophenol	12.746	196	377811	38.540	ng/ul	97
42) 2,4,5-Trichlorophenol	12.816	196	429220	38.769	ng/ul	99
43) 1,1'-Biphenyl	13.151	154	1458808	35.208	ng/ul	98
44) 2-Chloronaphthalene	13.193	162	1129058	34.616	ng/ul	99
45) 2-Nitroaniline	13.404	65	342072	39.408	ng/ul	98
47) Dimethylphthalate	13.787	163	1448039	34.539	ng/ul	100
48) 2,6-Dinitrotoluene	13.904	165	301435	38.779	ng/ul	98
50) Acenaphthylene	14.046	152	1858647	34.472	ng/ul	99
51) 3-Nitroaniline	14.234	138	347514	42.120	ng/ul	97
52) Acenaphthene	14.387	153	1207420	33.604	ng/ul	97
53) 2,4-Dinitrophenol	14.440	184	139971	39.204	ng/ul	94
55) 4-Nitrophenol	14.551	109	231755	37.680	ng/ul	98
56) Dibenzofuran	14.722	168	1734953	34.737	ng/ul	100
57) 2,4-Dinitrotoluene	14.693	165	458006	40.358	ng/ul	98
58) 2,3,4,6-Tetrachlorophenol	14.946	232	368285	40.054	ng/ul	96
59) Diethylphthalate	15.157	149	1498942	35.193	ng/ul	99
61) Fluorene	15.375	166	1378846	34.732	ng/ul	100
62) 4-Chlorophenyl-phenyle...	15.369	204	653138	33.534	ng/ul	95
63) 4-Nitroaniline	15.404	138	366922	45.744	ng/ul	98
66) 4,6-Dinitro-2-methylph...	15.457	198	246154	39.631	ng/ul	94
67) N-Nitrosodiphenylamine	15.587	169	1293009	36.254	ng/ul	98
68) 4-Bromophenyl-phenylether	16.269	248	431037	34.661	ng/ul	99
69) Hexachlorobenzene	16.369	284	505221	34.502	ng/ul	97
70) Atrazine	16.545	200	445287	36.021	ng/ul	98
71) Pentachlorophenol	16.716	266	333989	38.684	ng/ul	96
72) Phenanthrene	17.116	178	2378034	35.343	ng/ul	99
74) Anthracene	17.204	178	2323836	34.165	ng/ul	97
75) 1,2,3,4-Tetrachloroben...	13.110	216	561800	34.918	ng/uL	99
76) Pentachlorobenzene	14.634	250	554342	33.921	ng/uL	96
77) Carbazole	17.481	167	2317325	38.849	ng/ul	98
78) Di-n-butylphthalate	18.045	149	2770950	38.079	ng/ul	99
80) Fluoranthene	19.134	202	2858748	34.806	ng/ul	98
82) Pyrene	19.498	202	3073099	35.428	ng/ul	99
83) Butylbenzylphthalate	20.410	149	1350483	40.769	ng/ul	96
84) 3,3'-Dichlorobenzidine	21.222	252	906480	38.389	ng/ul	99
85) Benzo(a)anthracene	21.292	228	3058672	35.663	ng/ul	98
86) Bis(2-ethylhexyl)phtha...	21.216	149	1880002	38.485	ng/ul	99
87) Chrysene	21.351	228	2862871	35.365	ng/ul	100
89) Di-n-octyl phthalate	22.333	149	3543938	40.281	ng/ul	100
90) Benzo(b)fluoranthene	23.316	252	3050957	37.315	ng/ul	99
91) Benzo(k)fluoranthene	23.374	252	2982665	35.921	ng/ul	99
93) Benzo(a)pyrene	24.098	252	2552357	33.046	ng/ul	99
94) Indeno(1,2,3-cd)pyrene	27.539	276	3175101m	38.029	ng/ul	
95) Dibenzo(a,h)anthracene	27.598	278	2632457	37.453	ng/ul	99
96) Benzo(g,h,i)perylene	28.562	276	2405510	36.700	ng/ul	98

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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BNA_N

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