

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
 Data File : BG056109.D
 Acq On : 25 Dec 2022 4: 37
 Operator : CG/JU
 Sample : N6017-10MSD
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :

BNA_G

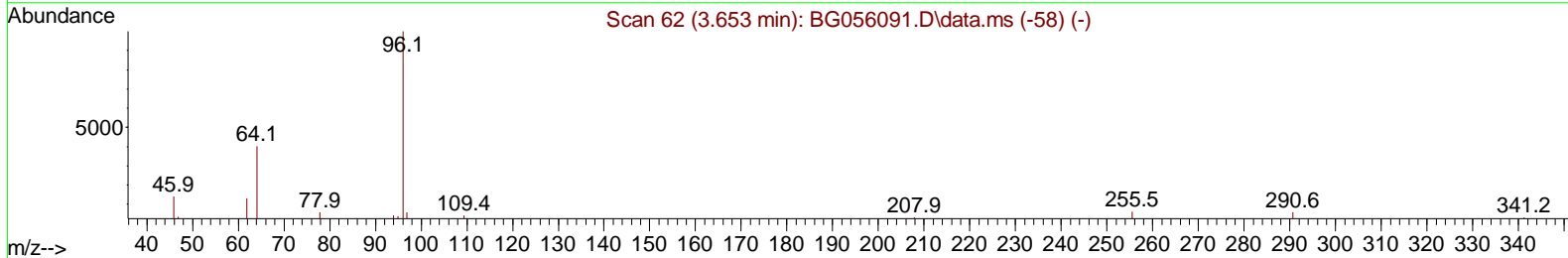
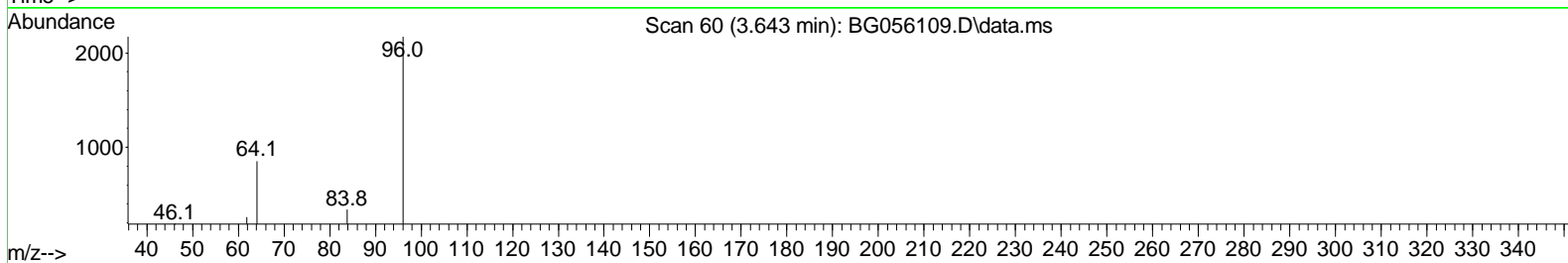
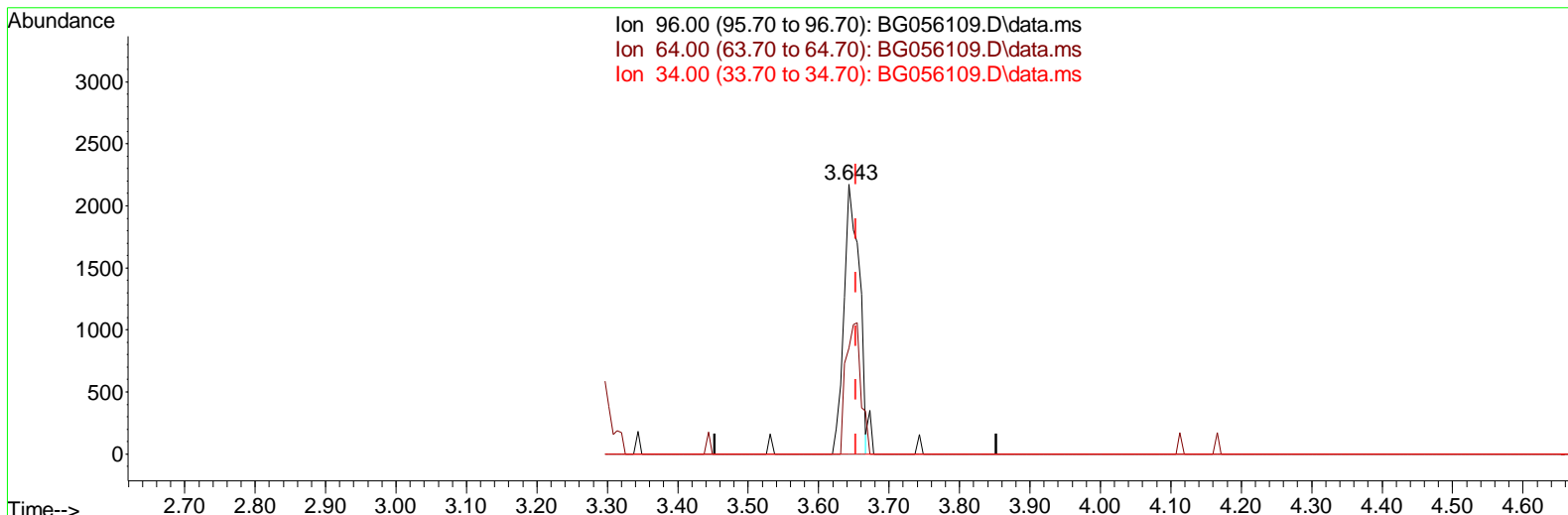
ClientSampleId :

DBYF6MSD

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/25/2022
 Supervised By :Yogesh Patel 12/25/2022

Quant Time: Dec 25 05: 51: 59 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG122422. M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Dec 24 22: 54: 38 2022
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(3) 1,4-Dioxane-d8 (S)

3.643min (-0.010) 3.26 ng/uL

response	3222	
Ion	Exp%	Act%
96.00	100.00	100.00
64.00	40.20	39.36
34.00	0.00	0.00
0.00	0.00	0.00

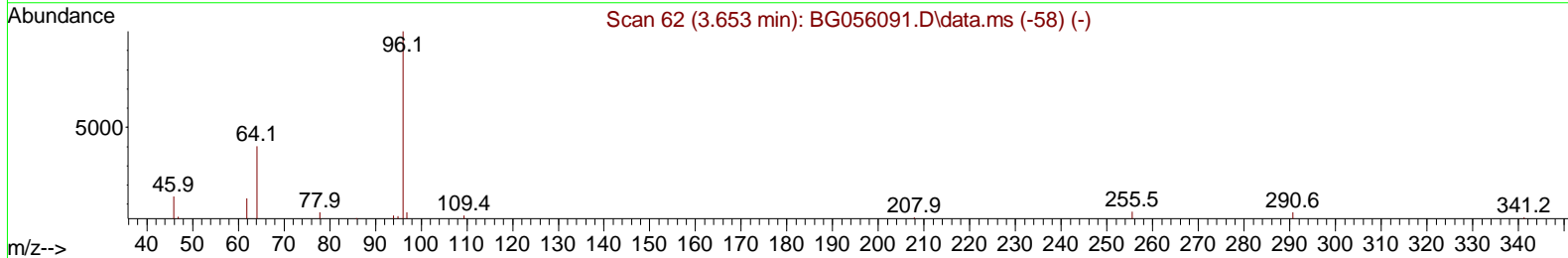
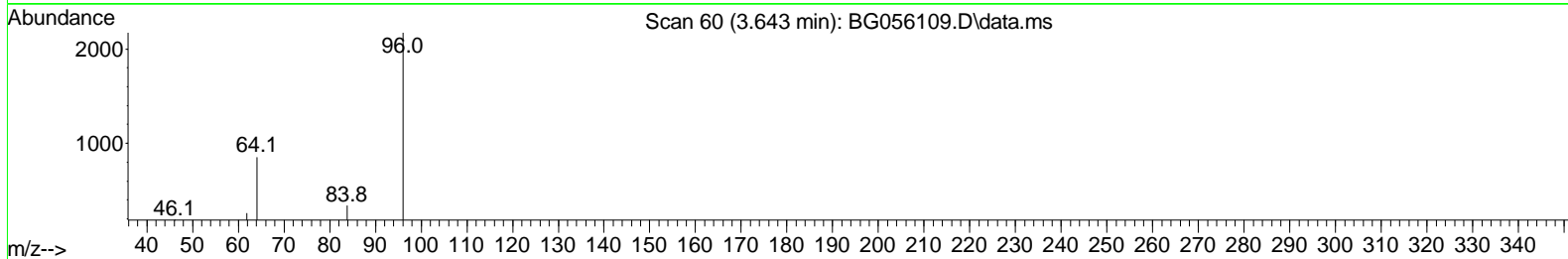
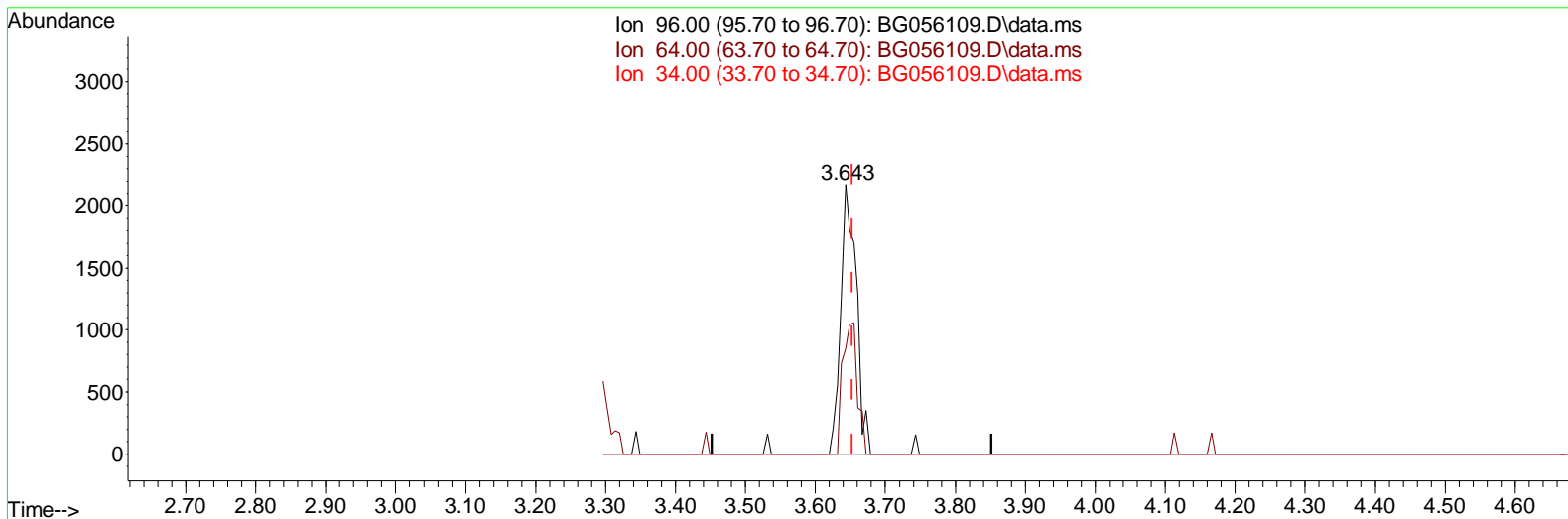
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(3) 1,4-Dioxane-d8 (S)

3.643min (-0.010) 3.39 ng/uL m

response	3346
Ion	Exp% Act%
96.00	100.00 100.00
64.00	40.20 39.36
34.00	0.00 0.00
0.00	0.00 0.00

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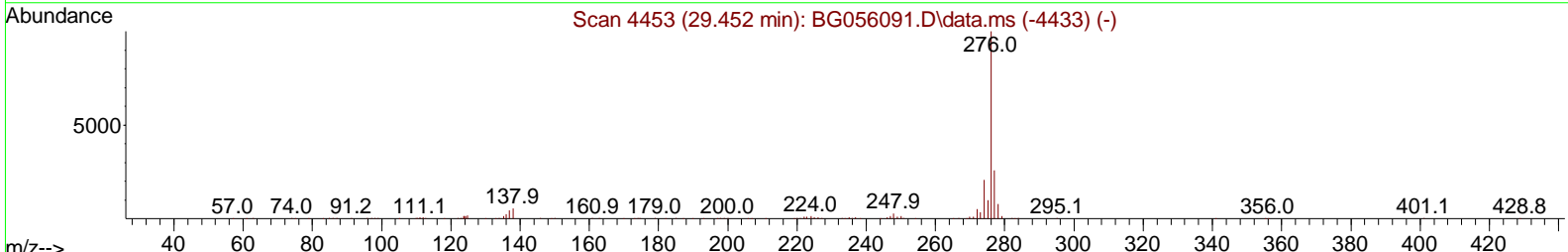
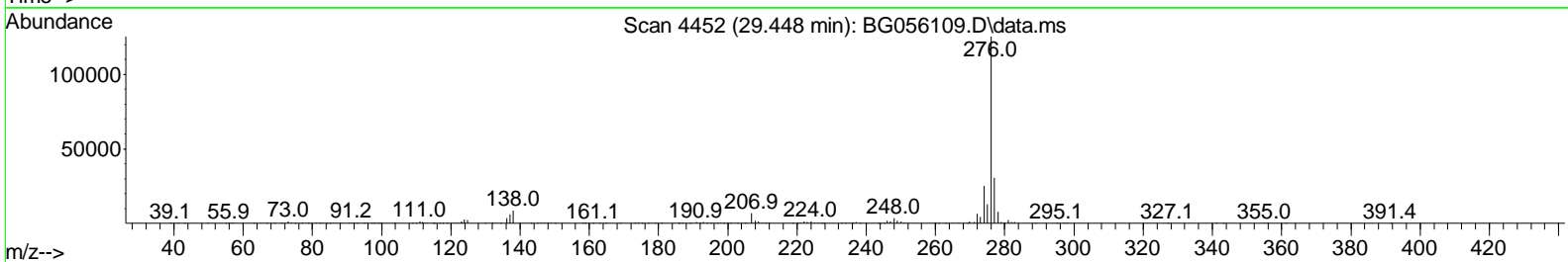
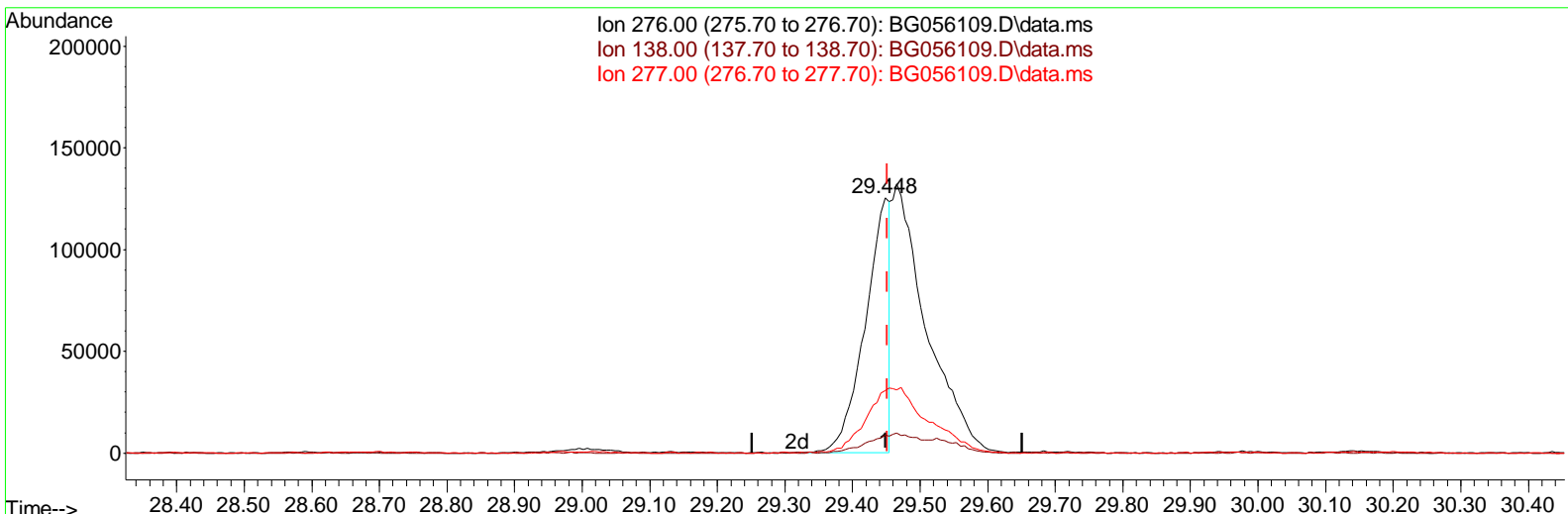
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(94) Indeno(1,2,3-cd)pyrene

29.448min (-0.004) 10.73 ng/ul

response 315242

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	5.50	7.02#
277.00	25.90	24.51
0.00	0.00	0.00

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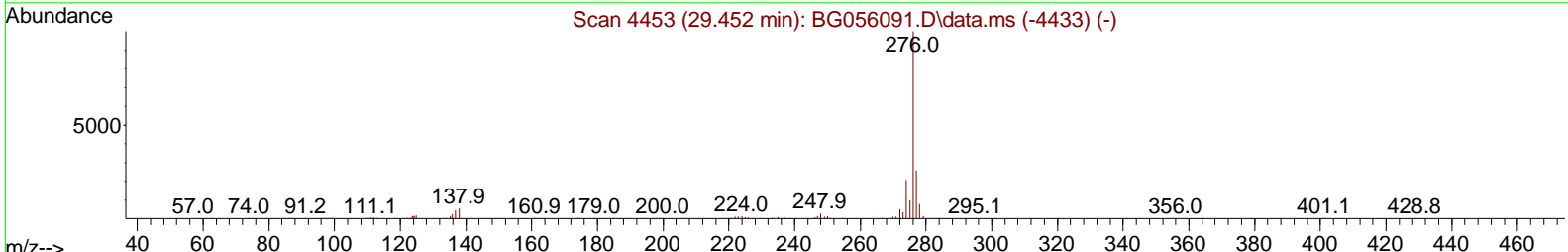
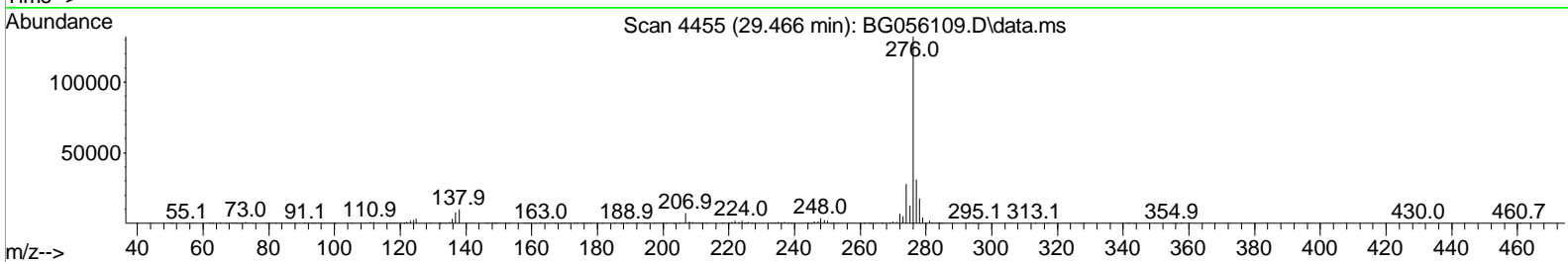
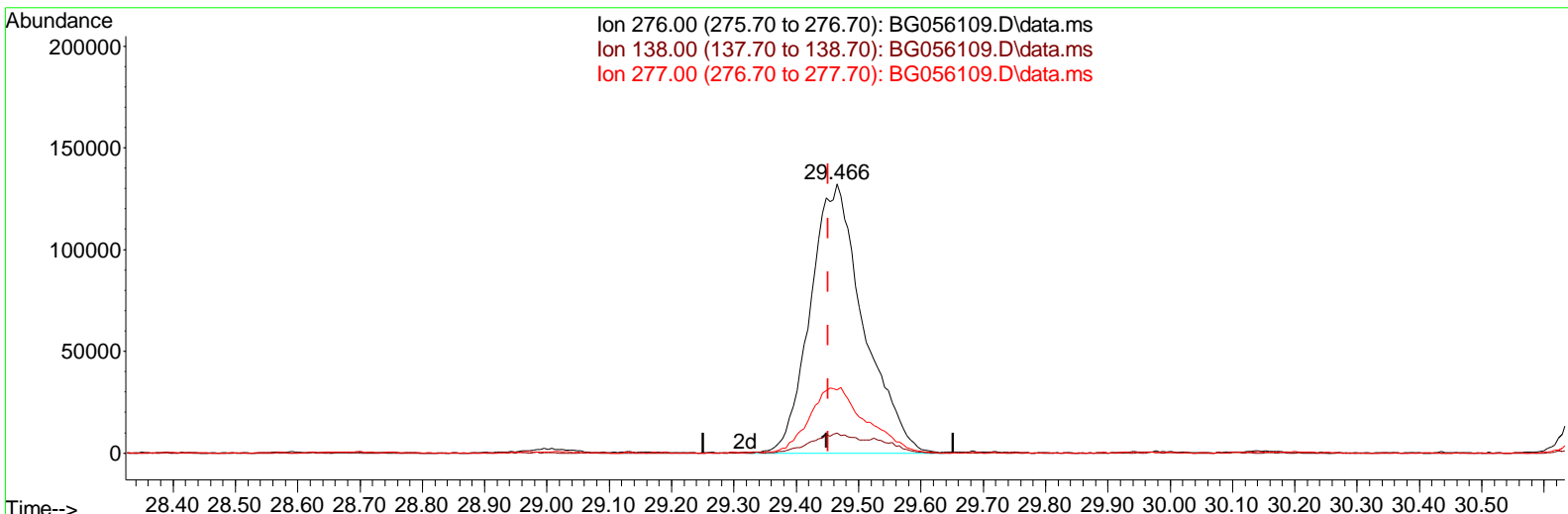
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(94) Indeno(1,2,3-cd)pyrene

29.466min (+ 0.014) 26.66 ng/ul m

response 783295

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	5.50	7.56#
277.00	25.90	23.48
0.00	0.00	0.00

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Compound	R. T.	QI on	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Di chlorobenzene-d4	8.343	152	41369	20.000	ng/ul	0.00
20) Naphthalene-d8	11.175	136	166671	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.953	164	144518	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.691	188	363448	20.000	ng/ul	0.00
79) Chrysene-d12	21.992	240	357857	20.000	ng/ul	0.00
88) Perylene-d12	25.464	264	387816	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.643	96	3346m	3.386	ng/uL	-0.01
4) Pyridine-d5	4.090	84	26734	8.297	ng/ul	0.00
7) Phenol-d5	7.474	99	77868	20.751	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.650	67	31452	21.266	ng/ul	0.00
11) 2-Chlorophenol-d4	7.867	132	51587	23.260	ng/ul	0.00
15) 4-Methylphenol-d8	9.037	113	56412	19.975	ng/ul	0.00
21) Nitrobenzene-d5	9.518	128	27803	22.903	ng/ul	0.00
24) 2-Nitrophenol-d4	10.247	143	35720	22.521	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.782	165	77798	22.074	ng/ul	0.00
31) 4-Chloroaniline-d4	11.305	131	68065	17.074	ng/ul	0.00
46) Dimethylphthalate-d6	14.342	166	253327	22.649	ng/ul	0.00
49) Acenaphthylene-d8	14.654	160	281479	22.284	ng/ul	0.00
54) 4-Nitrophenol-d4	15.118	143	37800	22.451	ng/ul	0.00
60) Fluorene-d10	15.934	176	248763	22.532	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.040	200	55151	23.605	ng/ul	0.00
73) Anthracene-d10	17.791	188	382668	22.784	ng/ul	0.00
81) Pyrene-d10	20.053	212	489353	24.760	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.224	264	465674	23.624	ng/ul	0.00
Target Compounds						
2) 1,4-Dioxane	3.690	88	7723	6.986	ng/uL#	92
5) Pyridine	4.107	79	33465	10.799	ng/ul #	94
6) Benzaldehyde	7.474	77	48451	35.169	ng/ul	88
8) Phenol	7.503	94	87262	23.366	ng/ul	98
10) Bis(2-Chloroethyl)ether	7.750	93	53852	21.906	ng/ul	98
12) 2-Chlorophenol	7.903	128	53592	24.057	ng/ul	97
13) 2-Methylphenol	8.772	108	63243	22.554	ng/ul	85
14) 2,2'-oxybis(1-Chloropr...	8.860	45	8382	26.546	ng/ul #	74
16) Acetophenone	9.172	105	115144	24.613	ng/ul	94
17) N-Nitrosodipropylamine	9.142	70	46024	23.269	ng/ul	96
18) 4-Methylphenol	9.101	108	71331	23.484	ng/ul	92
19) Hexachloroethane	9.448	117	20694	23.104	ng/ul	98
22) Nitrobenzene	9.560	77	68583	22.768	ng/ul	95
23) Isophorone	10.082	82	147386	21.400	ng/ul	98
25) 2-Nitrophenol	10.276	139	36751	24.004	ng/ul #	84
26) 2,4-Dimethylphenol	10.318	107	77017	21.259	ng/ul	98
27) Bis(2-Chloroethoxy)meth...	10.558	93	77943	21.838	ng/ul #	96
29) 2,4-Dichlorophenol	10.811	162	77156	22.921	ng/ul #	86
30) Naphthalene	11.228	128	198596	22.868	ng/ul	99
32) 4-Chloroaniline	11.328	127	76149	18.956	ng/ul	96
33) Hexachlorobutadiene	11.510	225	64807	20.544	ng/ul	96
34) Caprolactam	12.068	113	24198	22.484	ng/ul #	79
35) 4-Chloro-3-methylphenol	12.415	107	79861	24.695	ng/ul	99

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Compound	R. T.	QI on	Response	Conc	Units	Dev(Mi n)
36) 2-Methyl naphthal ene	12. 809	142	156950	22. 542	ng/ul	100
37) 1-Methyl naphthal ene	13. 020	142	154832	22. 075	ng/ul	95
39) 1, 2, 4, 5-Tetrachl oroben. . .	13. 167	216	132278	22. 194	ng/ul	97
40) Hexachl orocycl opentadi ene	13. 144	237	63109	16. 331	ng/ul	99
41) 2, 4, 6-Tri chl orophenol	13. 396	196	84064	23. 930	ng/ul	93
42) 2, 4, 5-Tri chl orophenol	13. 467	196	91328	23. 978	ng/ul	90
43) 1, 1' -Bi phenyl	13. 790	154	232942	22. 800	ng/ul	98
44) 2-Chl oronaphthal ene	13. 843	162	195246	23. 056	ng/ul	97
45) 2-Ni troani li ne	14. 031	65	35088	29. 837	ng/ul	79
47) Di methyl phthal ate	14. 389	163	258252	23. 284	ng/ul	100
48) 2, 6-Di ni trotol uene	14. 513	165	53670	24. 515	ng/ul	94
50) Acenaphthyl ene	14. 683	152	294200	23. 437	ng/ul	98
51) 3-Ni troani li ne	14. 842	138	44342	26. 965	ng/ul	100
52) Acenaphthene	15. 018	153	208193	23. 609	ng/ul	94
53) 2, 4-Di ni trophenol	15. 047	184	36012	21. 134	ng/ul #	90
55) 4-Ni trophenol	15. 130	109	40272	24. 586	ng/ul	93
56) Di benzofuran	15. 347	168	311914	22. 597	ng/ul	99
57) 2, 4-Di ni trotol uene	15. 294	165	75258	24. 789	ng/ul #	96
58) 2, 3, 4, 6-Tetrachl orophenol	15. 564	232	88154	23. 414	ng/ul #	98
59) Di ethyl phthal ate	15. 741	149	235945	23. 811	ng/ul	100
61) Fl uorene	15. 993	166	257162	22. 680	ng/ul	98
62) 4-Chl orophenyl -phenyl e. . .	15. 976	204	159788	21. 693	ng/ul	96
63) 4-Ni troani li ne	15. 999	138	44907	28. 619	ng/ul #	77
66) 4, 6-Di ni tro-2-methyl ph. . .	16. 058	198	54621	24. 384	ng/ul #	91
67) N-Ni trosodi phenyl ami ne	16. 187	169	233549	25. 362	ng/ul	96
68) 4-Bromophenyl -phenyl ether	16. 869	248	110090	23. 892	ng/ul	95
69) Hexachl orobenzene	16. 998	284	108649	24. 363	ng/ul	95
70) Atrazi ne	17. 121	200	108456	24. 701	ng/ul	96
71) Pentachl orophenol	17. 333	266	58865	20. 761	ng/ul	91
72) Phenanthrene	17. 732	178	486460	26. 365	ng/ul	100
74) Anthracene	17. 826	178	445513	23. 587	ng/ul	98
75) 1, 2, 3, 4-Tetrachl oroben. . .	13. 766	216	136917	24. 836	ng/uL	97
76) Pentachl orobenzene	15. 271	250	131575	24. 177	ng/uL	91
77) Carbazol e	18. 085	167	383557	24. 838	ng/ul	98
78) Di -n-butyl phthal ate	18. 620	149	393906	27. 919	ng/ul	99
80) Fl uoranthene	19. 724	202	728813	31. 361	ng/ul	99
82) Pyrene	20. 083	202	685883	29. 256	ng/ul #	97
83) Butyl benzyl phthal ate	20. 946	149	166403	33. 354	ng/ul #	88
84) 3, 3' -Di chl orobenzi di ne	21. 869	252	147729	18. 373	ng/ul	99
85) Benzo(a)anthracene	21. 969	228	687479	28. 146	ng/ul	99
86) Bi s(2-ethyl hexyl)phtha. . .	21. 839	149	242583	31. 740	ng/ul	99
87) Chrysene	22. 045	228	638359	28. 161	ng/ul	98
89) Di -n-octyl phthal ate	23. 138	149	416311	29. 008	ng/ul	100
90) Benzo(b)fl uoranthene	24. 354	252	744965	29. 098	ng/ul	99
91) Benzo(k)fl uoranthene	24. 424	252	661829	26. 491	ng/ul #	99
93) Benzo(a)pyrene	25. 300	252	620549	27. 816	ng/ul #	99
94) I ndeno(1, 2, 3-cd)pyrene	29. 466	276	783295m	26. 658	ng/ul	98
95) Di benzo(a, h)anthracene	29. 530	278	629249	25. 603	ng/ul	98
96) Benzo(g, h, i)peryl ene	30. 723	276	603126	25. 541	ng/ul	98

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

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