

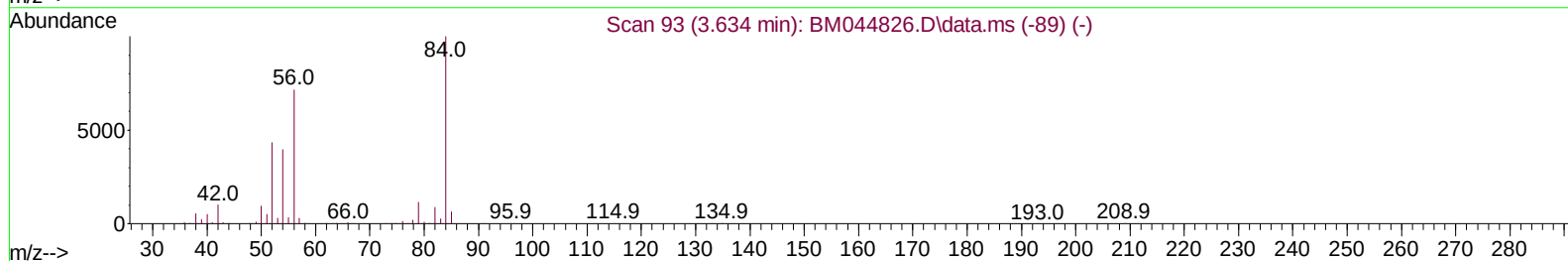
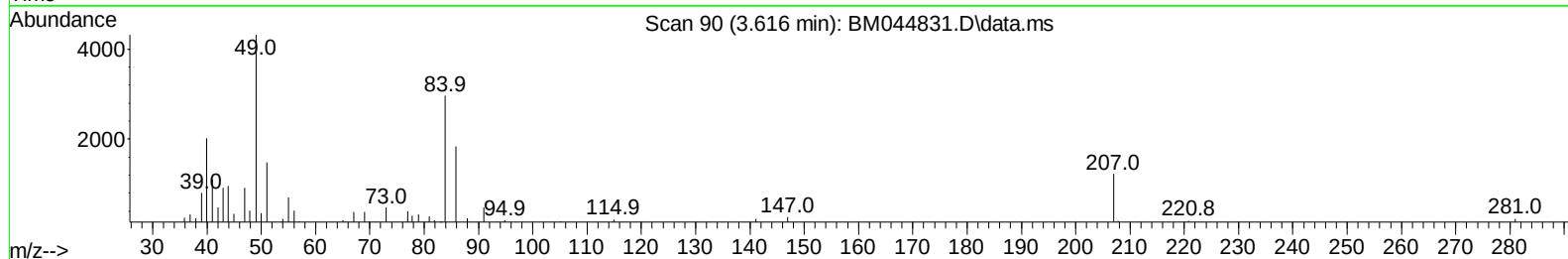
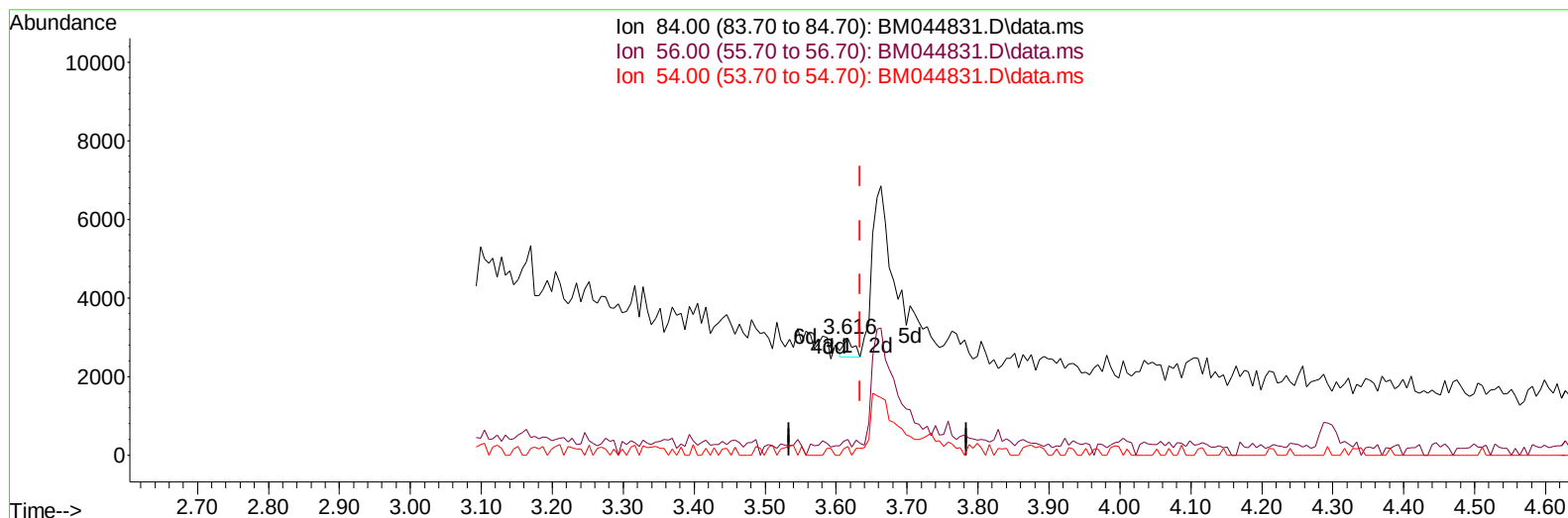
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM033024\
Data File : BM044831.D
Acq On : 30 Mar 2024 14:27
Operator : MA/JU
Sample : P1887-04
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
E0797

Manual IntegrationsAPPROVED

Quant Time: Mar 31 15:15:33 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030524.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sun Mar 31 13:54:26 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/01/2024
Supervised By :mohammad ahmed 04/02/2024



TIC: BM044831.D\data.ms

(4) Pyridine-d5 (S)

3.616min (-0.018) 0.06 ng/u1

response 471

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	63.10	13.65#
54.00	29.60	7.47#
0.00	0.00	0.00

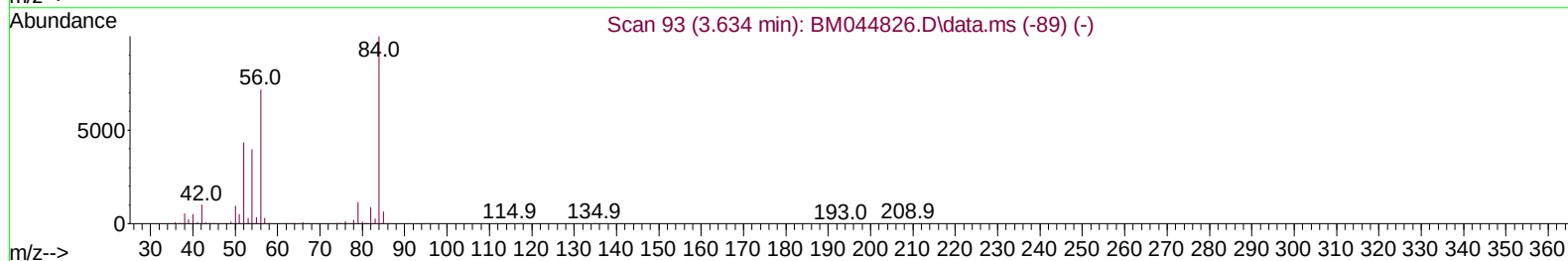
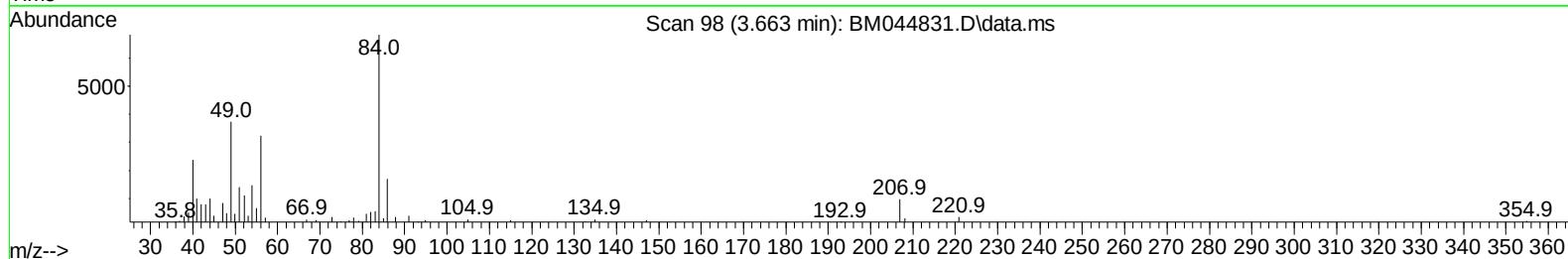
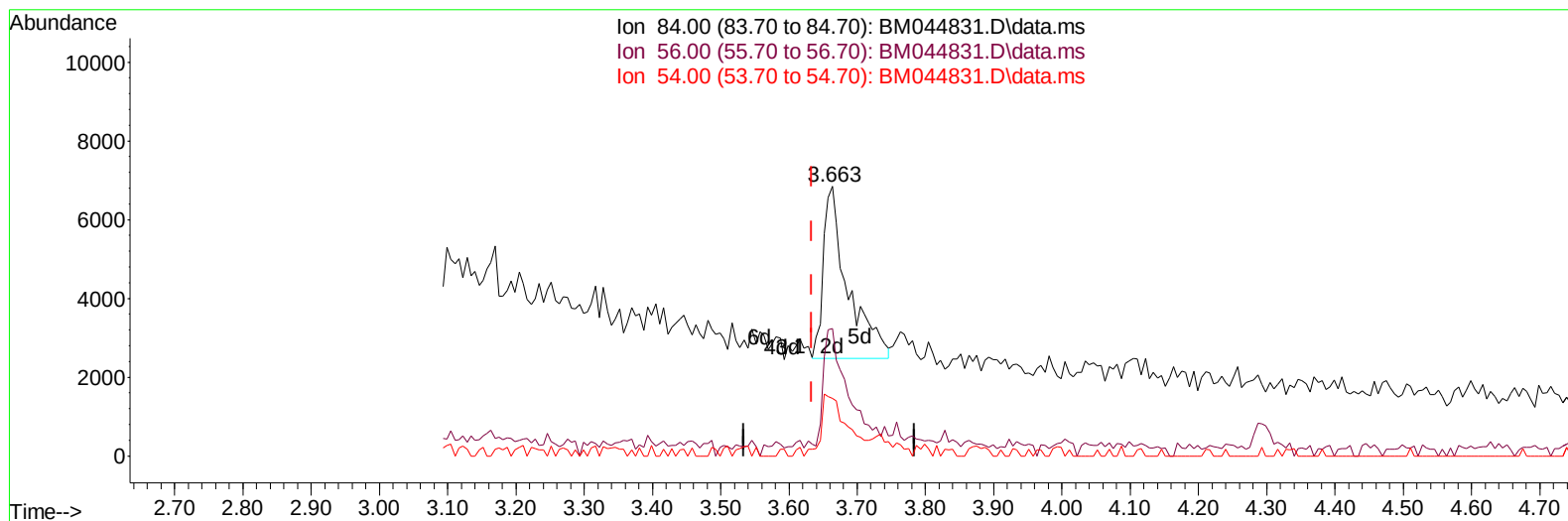
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(4) Pyridine-d5 (S)

3.663min (+ 0.030) 1.42 ng/ul m

response 10876

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	63.10	47.21#
54.00	29.60	21.34#
0.00	0.00	0.00

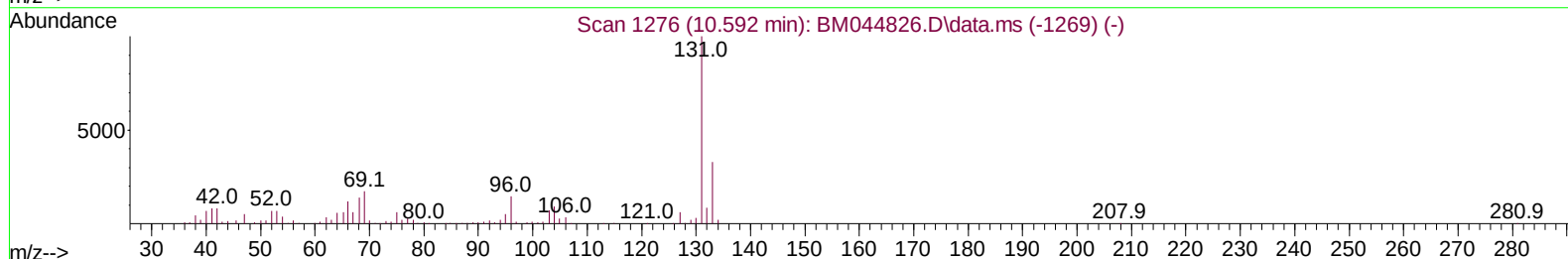
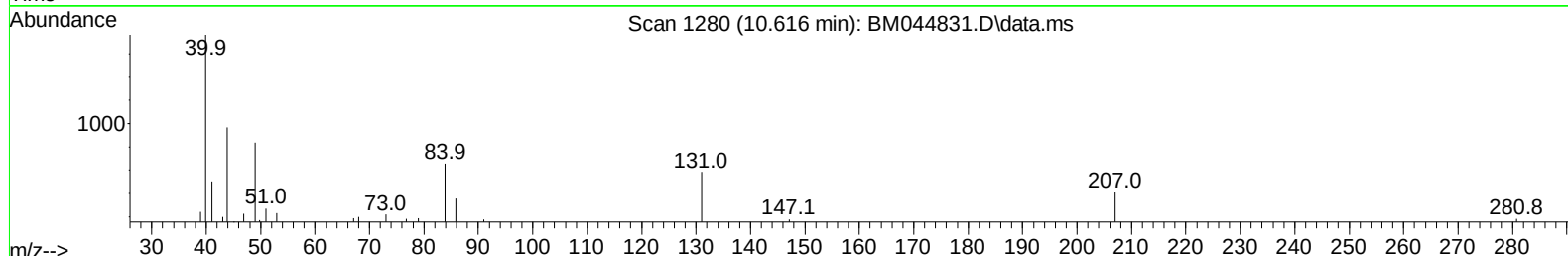
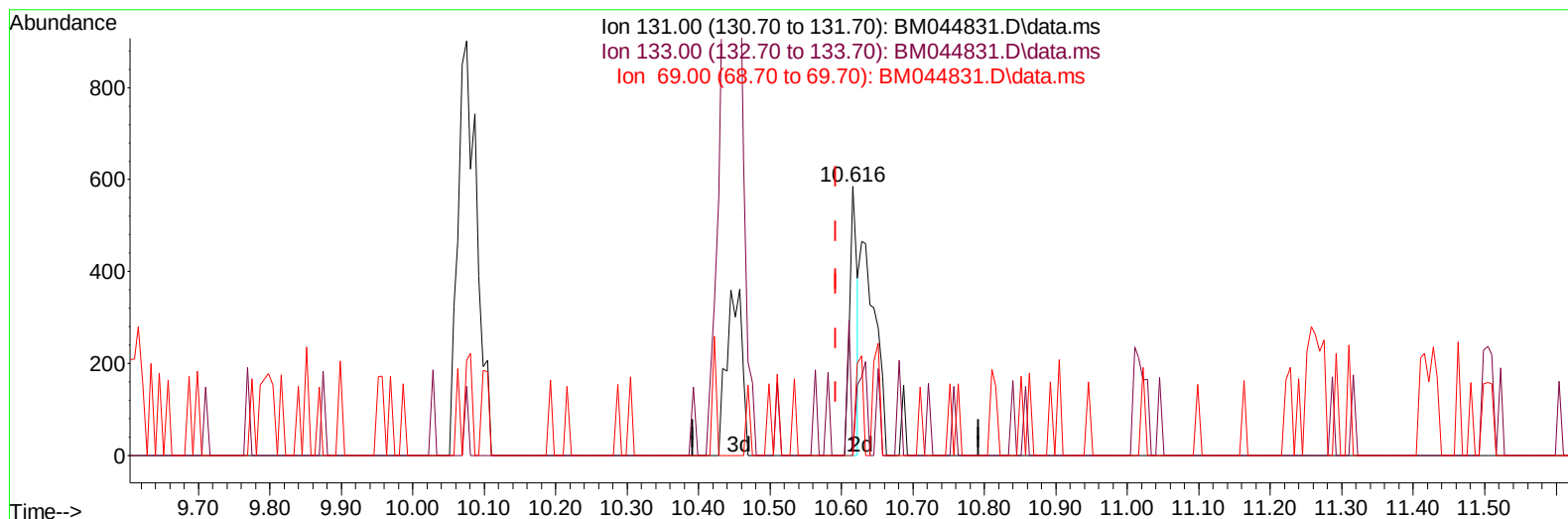
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ALS Vial : 7 Sample Multiplier: 1

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TIC: BM044831.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.616min (+ 0.024) 0.05 ng/uI

response 434

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.40	0.00#
69.00	15.90	0.00#
0.00	0.00	0.00

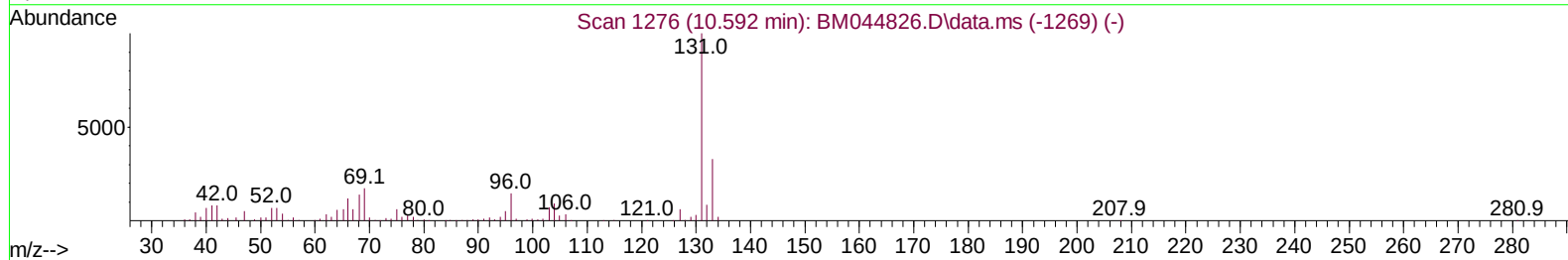
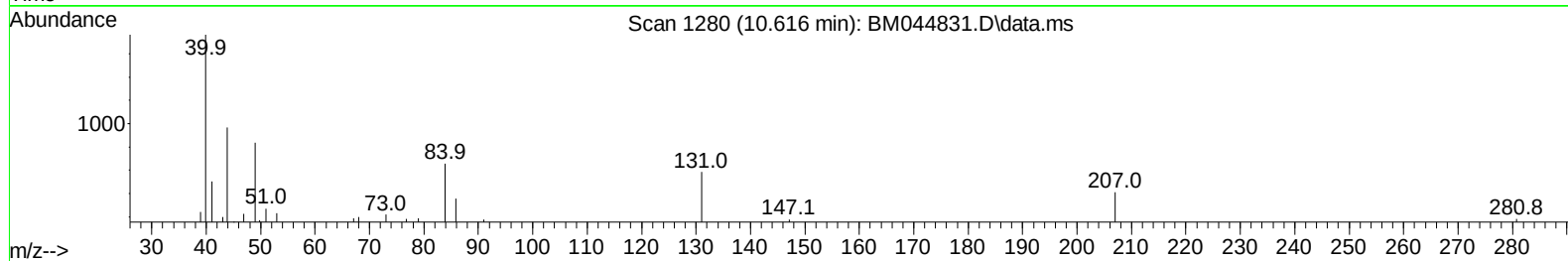
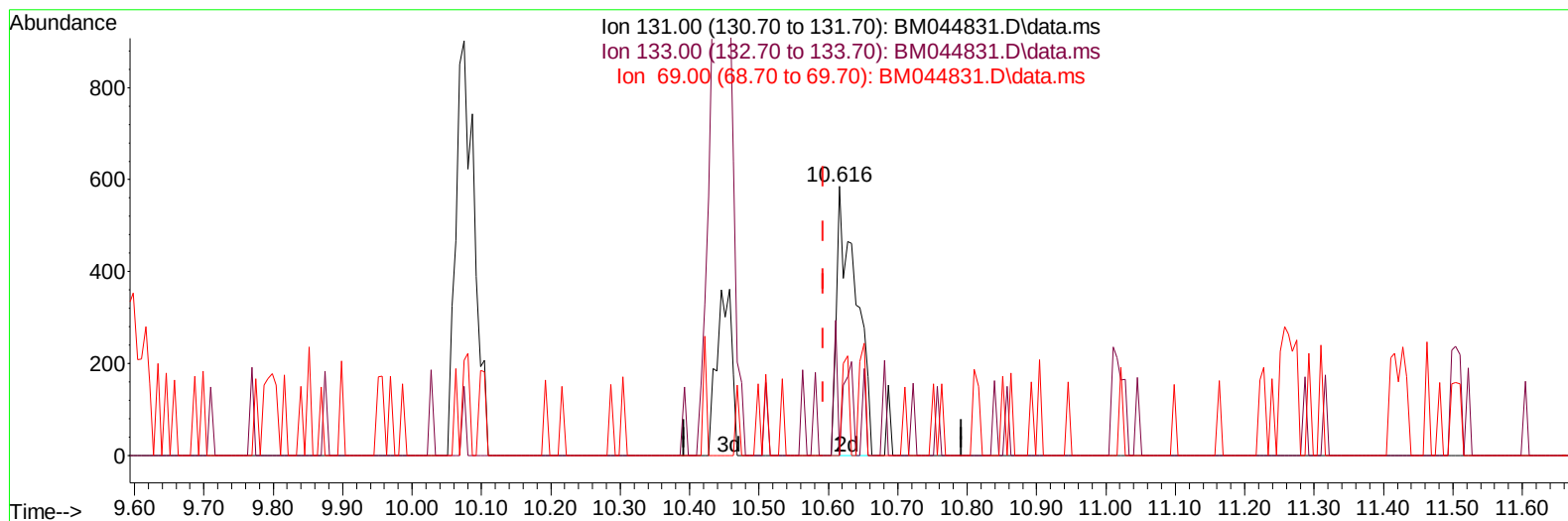
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TIC: BM044831.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.616min (+ 0.024) 0.13 ng/ul m

response 1149

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.40	0.00#
69.00	15.90	0.00#
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/01/2024
 Supervised By :mohammad ahmed 04/02/2024

Quant Time: Mar 31 15:16:46 2024
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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.669	152		87955	20.000	ng/ul	0.00
20) Naphthalene-d8	10.445	136		349470	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.310	164		199318	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.068	188		388080	20.000	ng/ul	0.00
79) Chrysene-d12	21.268	240		366790	20.000	ng/ul	0.00
88) Perylene-d12	23.503	264		415050	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.234	96		5922	2.284	ng/uL	0.00
4) Pyridine-d5	3.663	84		10876m	1.425	ng/ul	0.03
7) Phenol-d5	6.846	99		89636	9.912	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.016	67		63062	11.290	ng/ul	0.00
11) 2-Chlorophenol-d4	7.199	132		73241	11.022	ng/ul	0.00
15) 4-Methylphenol-d8	8.375	113		31597	4.449	ng/ul	0.00
21) Nitrobenzene-d5	8.828	128		37482	12.737	ng/ul	0.00
24) 2-Nitrophenol-d4	9.540	143		40226	12.680	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.075	165		59380	10.998	ng/ul	0.00
31) 4-Chloroaniline-d4	10.616	131		1149m	0.128	ng/ul	0.02
46) Dimethylphthalate-d6	13.733	166		198448	12.692	ng/ul	0.00
49) Acenaphthylene-d8	14.004	160		229696	12.846	ng/ul	0.00
54) 4-Nitrophenol-d4	14.551	143		25597	7.894	ng/ul	0.02
60) Fluorene-d10	15.310	176		169310	12.886	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.451	200		17142	6.805	ng/ul	0.00
73) Anthracene-d10	17.168	188		236777	12.764	ng/ul	0.00
81) Pyrene-d10	19.468	212		292004	12.980	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.362	264		288575	13.228	ng/ul	0.00
Target Compounds							
16) Acetophenone	8.493	105		23035	2.091	ng/ul#	Qvalue 93

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

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