

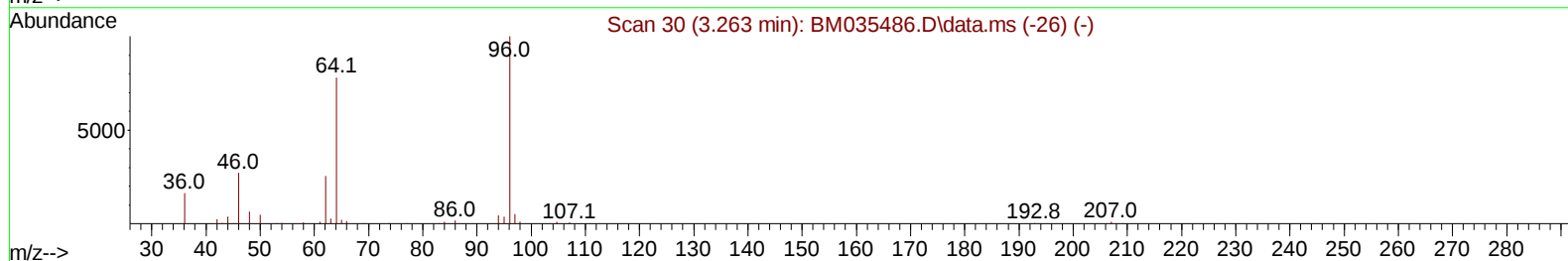
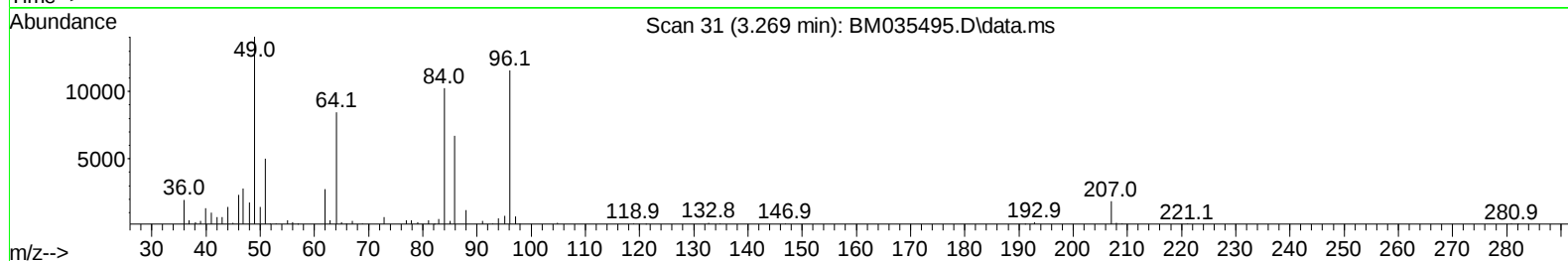
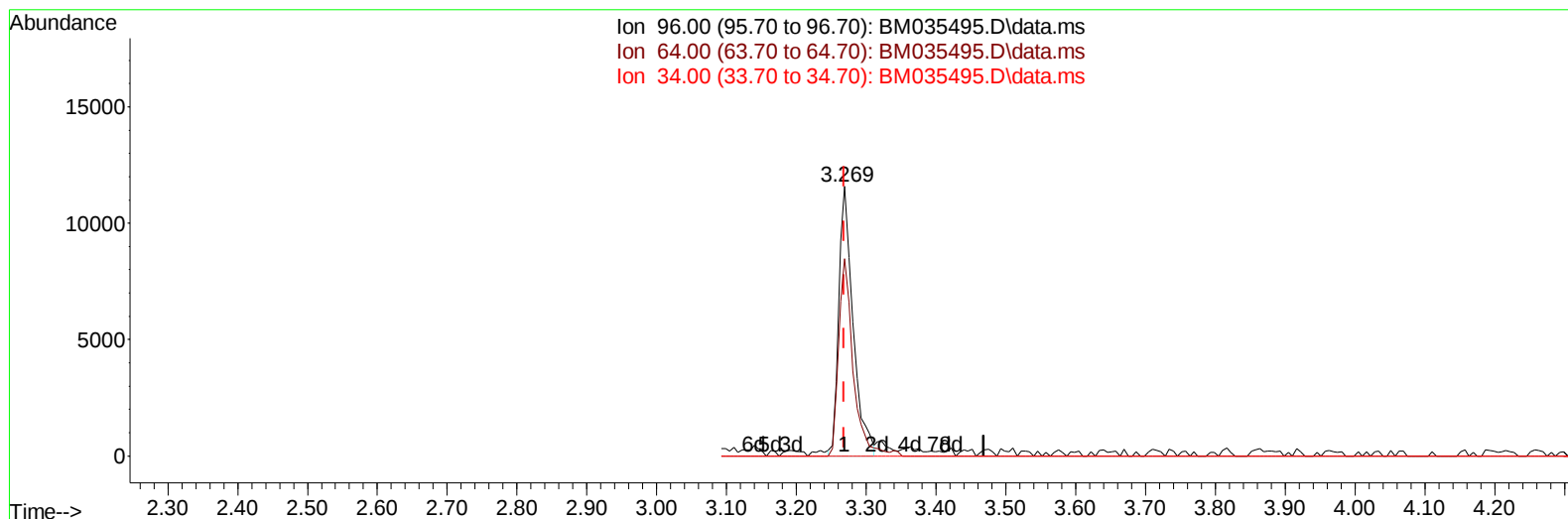
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061722\
Data File : BM035495.D
Acq On : 17 Jun 2022 15:12
Operator : CG/JU
Sample : N3226-14
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EW4N4

Manual IntegrationsAPPROVED

Quant Time: Jun 17 23:01:52 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM061622.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jun 17 01:58:49 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/18/2022
Supervised By :Yogesh Patel 06/22/2022



TIC: BM035495.D\data.ms

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

3.269min (+ 0.000) 3.37 ng/uL

response 16379

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	75.40	73.19
34.00	0.00	0.00
0.00	0.00	0.00

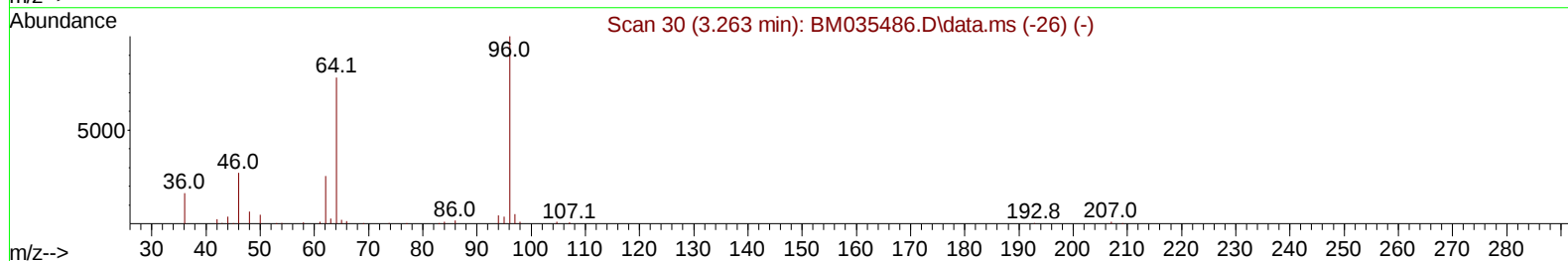
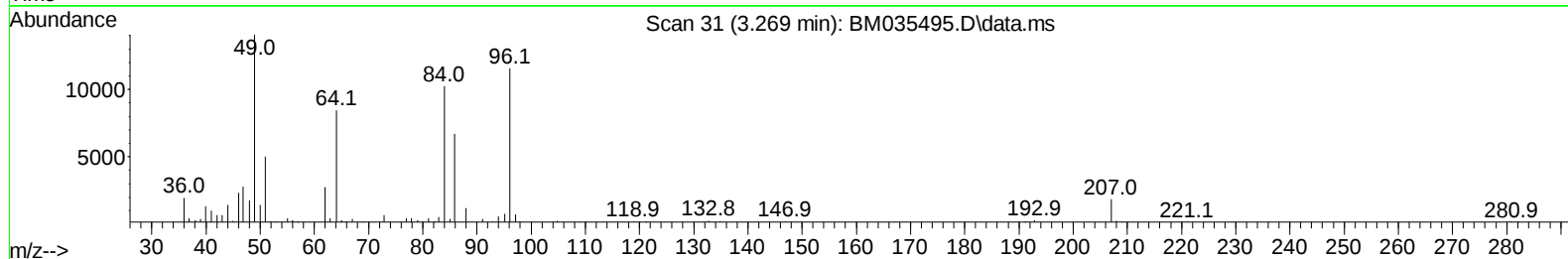
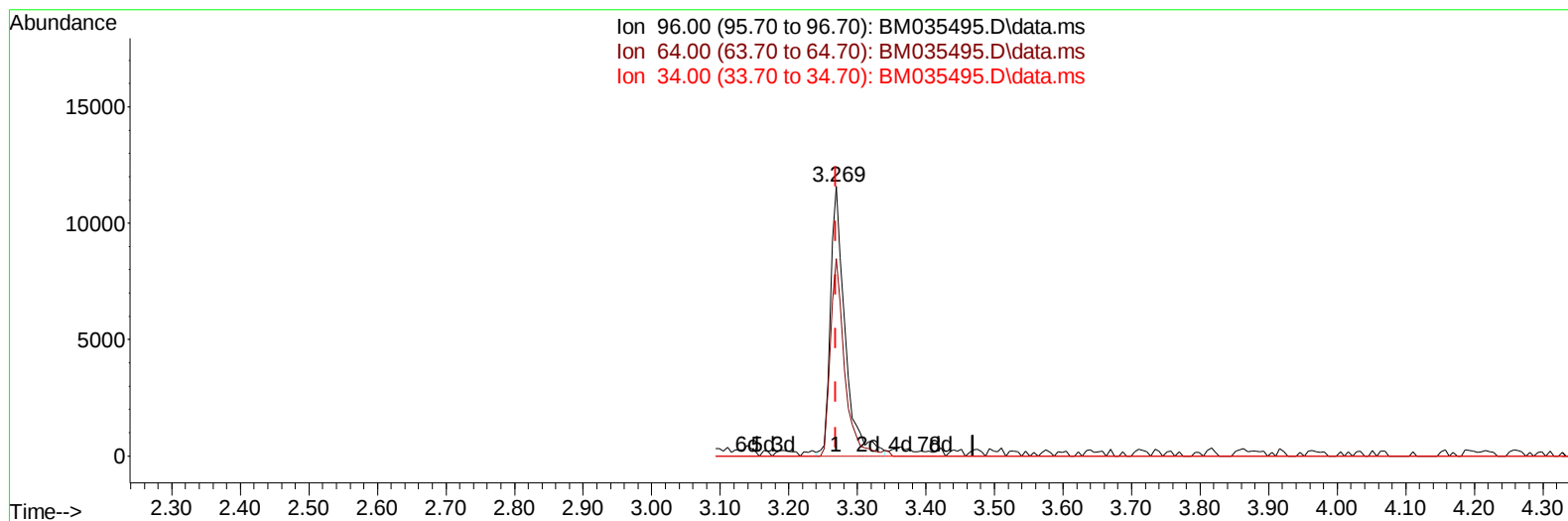
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TIC: BM035495.D\data.ms

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

3.269min (+ 0.000) 3.56 ng/uL m

response 17255

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	75.40	73.19
34.00	0.00	0.00
0.00	0.00	0.00

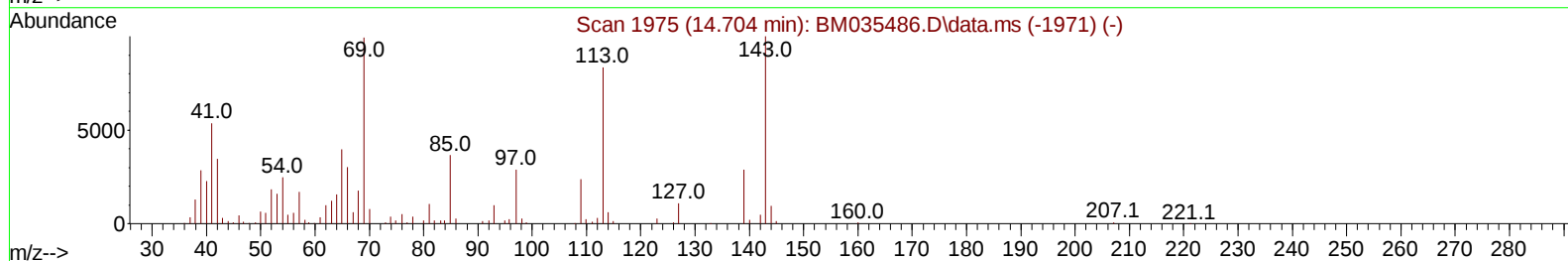
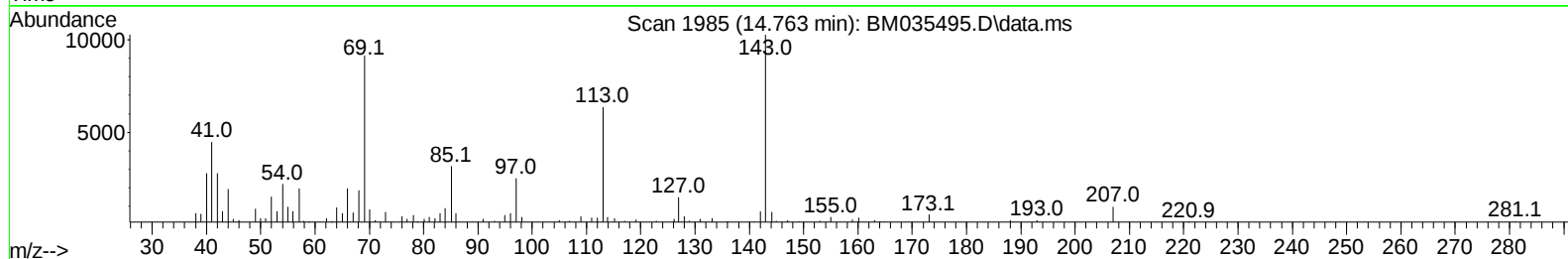
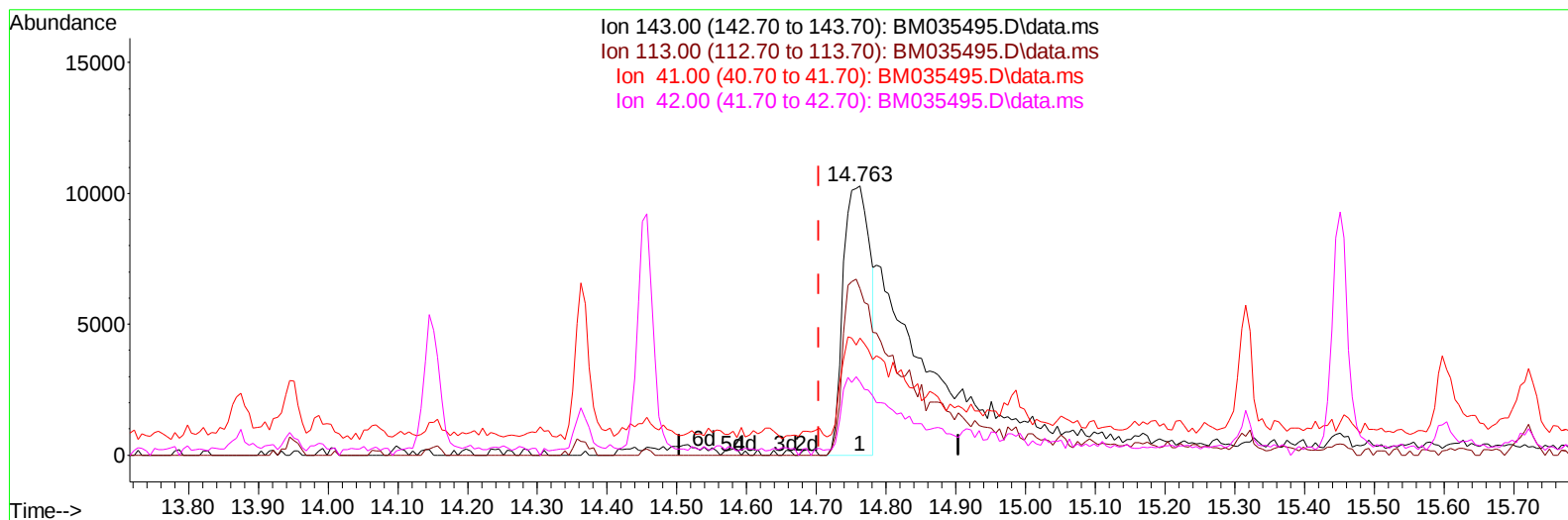
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TIC: BM035495.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (+ 0.059) 5.25 ng/u1

response 27131

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	61.40	61.91
41.00	41.00	43.59
42.00	29.30	27.20

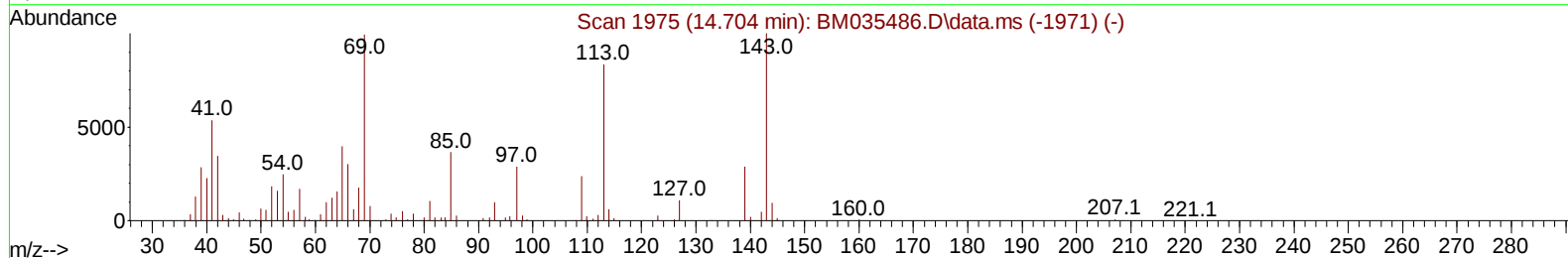
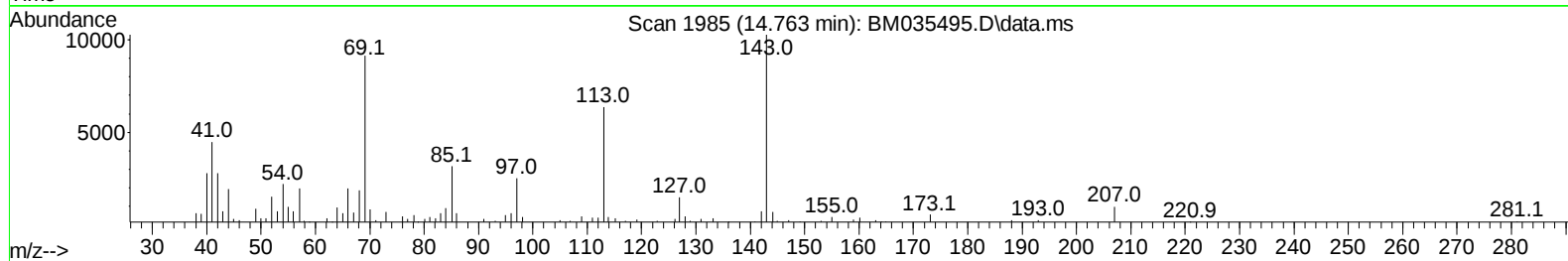
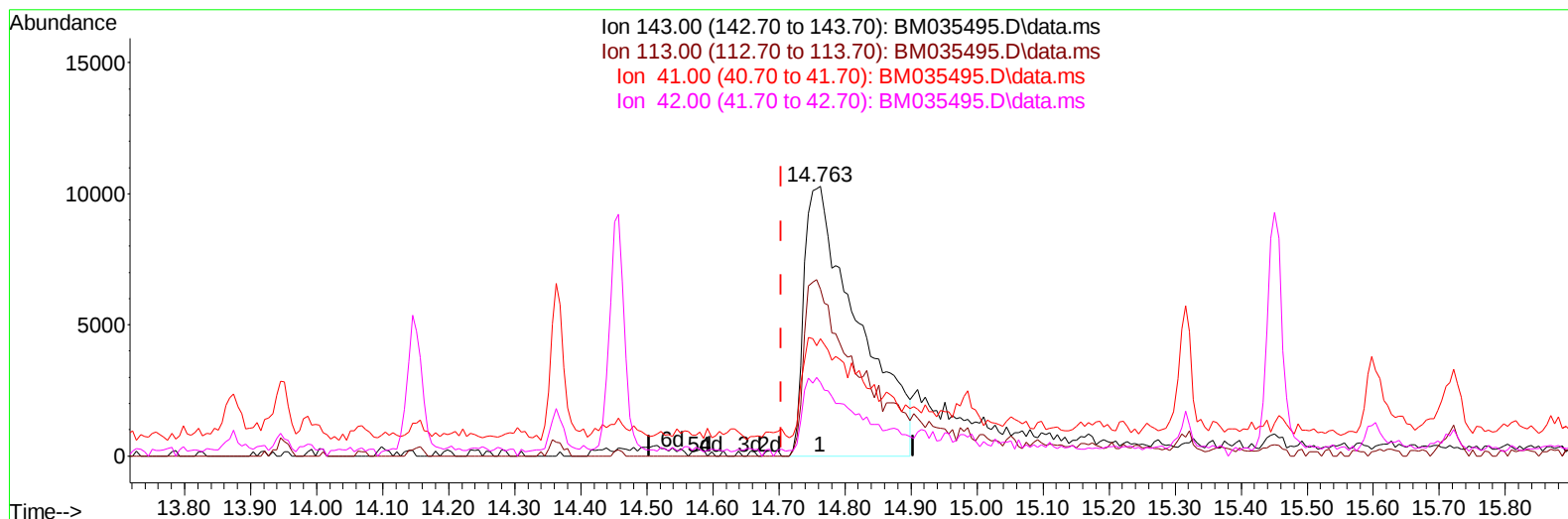
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Instrument :
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Manual IntegrationsAPPROVED

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TIC: BM035495.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (+ 0.059) 11.11 ng/ul m

response 57478

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	61.40	61.91
41.00	41.00	43.59
42.00	29.30	27.20

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 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Manual IntegrationsAPPROVED

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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.816	152		193744	20.000	ng/ul	0.00
20) Naphthalene-d8	10.610	136		818511	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.457	164		494944	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.204	188		1016536	20.000	ng/ul	0.00
79) Chrysene-d12	21.392	240		1023780	20.000	ng/ul	0.00
88) Perylene-d12	23.733	264		1080773	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.269	96		17255m	3.555	ng/uL	0.00
4) Pyridine-d5	3.699	84		97204	8.132	ng/ul	0.01
7) Phenol-d5	7.010	99		236327	15.916	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.151	67		159534	17.277	ng/ul	0.00
11) 2-Chlorophenol-d4	7.357	132		231313	17.986	ng/ul	0.00
15) 4-Methylphenol-d8	8.545	113		200645	16.142	ng/ul	0.00
21) Nitrobenzene-d5	8.981	128		104907	17.218	ng/ul	0.00
24) 2-Nitrophenol-d4	9.710	143		110128	16.762	ng/ul	0.01
28) 2,4-Dichlorophenol-d3	10.263	165		165918	14.874	ng/ul	0.01
31) 4-Chloroaniline-d4	10.769	131		246349	14.036	ng/ul	0.01
46) Dimethylphthalate-d6	13.874	166		639864	19.006	ng/ul	0.00
49) Acenaphthylene-d8	14.151	160		756609	18.135	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143		57478m	11.114	ng/ul	0.06
60) Fluorene-d10	15.451	176		573433	19.671	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.604	200		60814	10.035	ng/ul	0.01
73) Anthracene-d10	17.304	188		908502	19.883	ng/ul	0.00
81) Pyrene-d10	19.598	212		1129204	22.435	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.580	264		1168112	22.408	ng/ul	0.00
Target Compounds							
16) Acetophenone	8.657	105		39180	2.079	ng/ul	Qvalue 94

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

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