Quantitation Report (Qedit)

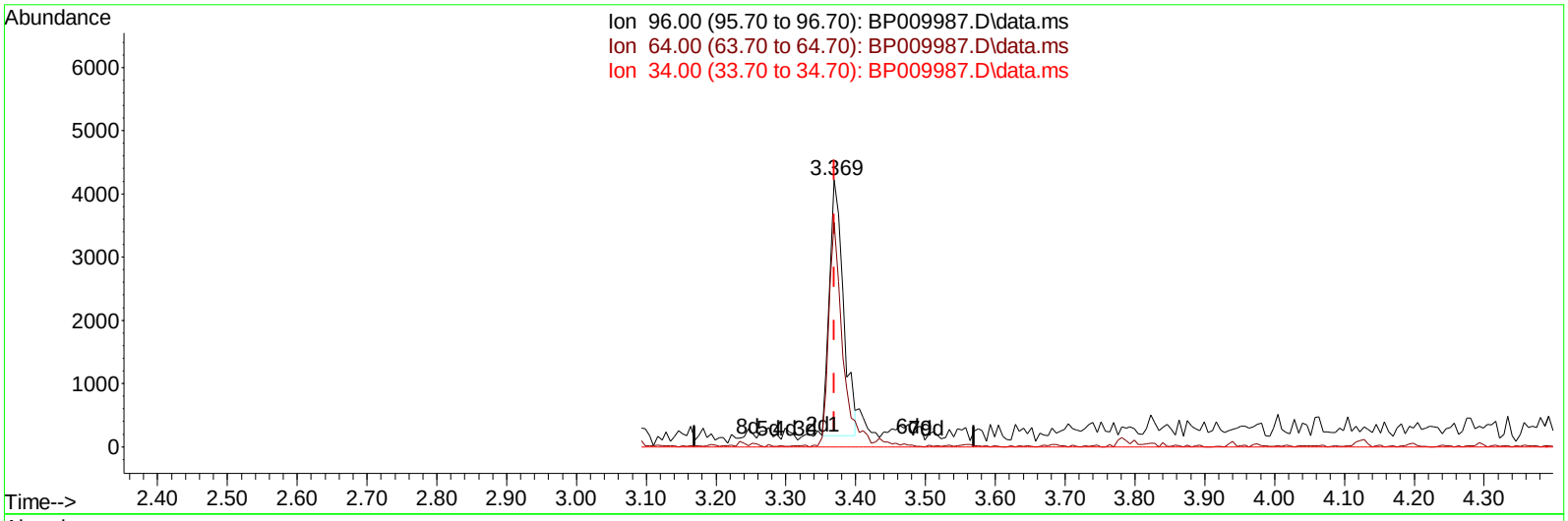
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
Data File : BP009987.D
Acq On : 20 Apr 2022 21:12
Operator : CG/JU
Sample : N2472-02
Misc :
ALS Vial : 40 Sample Multiplier: 1

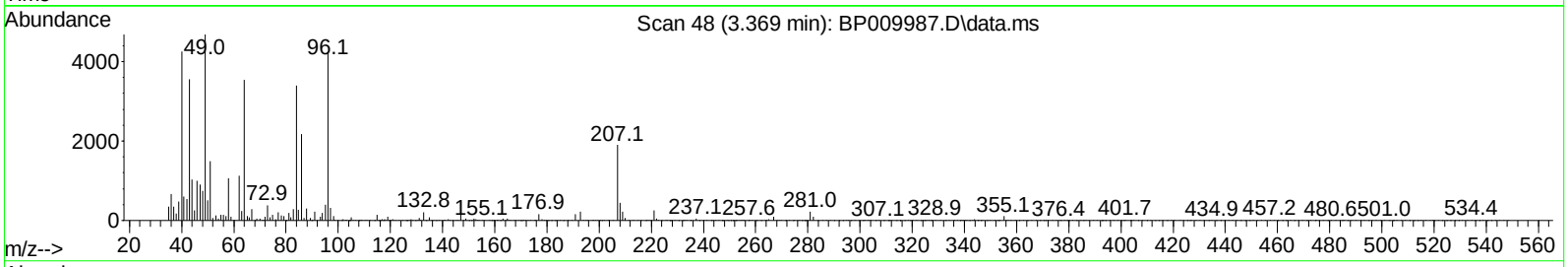
Quant Time: Apr 21 01:09:07 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Apr 20 01:32:14 2022
Response via : Initial Calibration

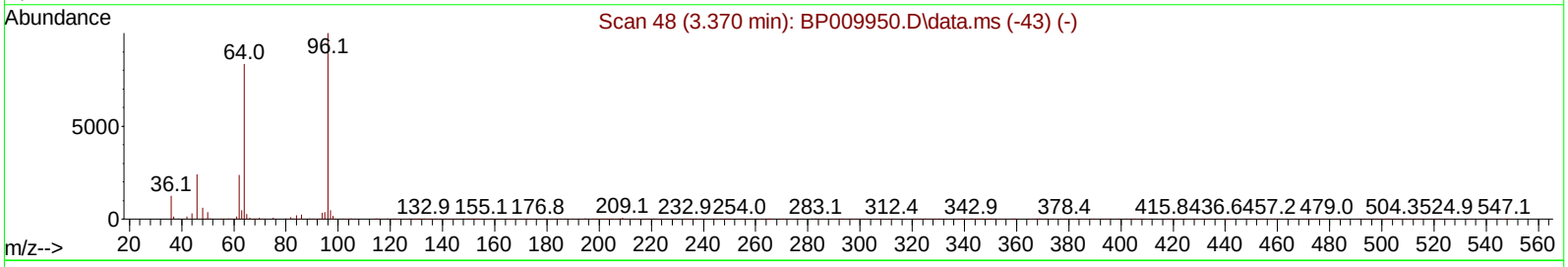
Instrument :
BNA_P
ClientSampleId :
C0J29

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/21/2022
Supervised By :Yogesh Patel 04/25/2022







TIC: BP009987.D\data.ms

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

3.369min (-0.000) 2.01 ng/uL

response 5586

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	84.04#
34.00	0.00	0.00
0.00	0.00	0.00

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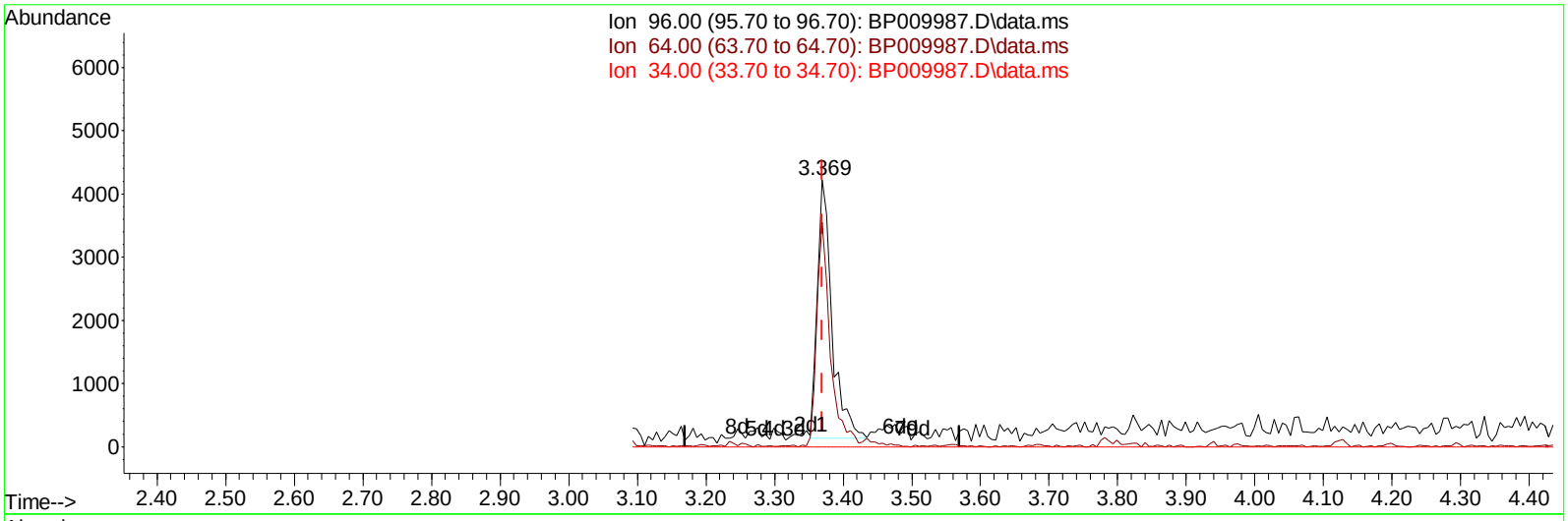

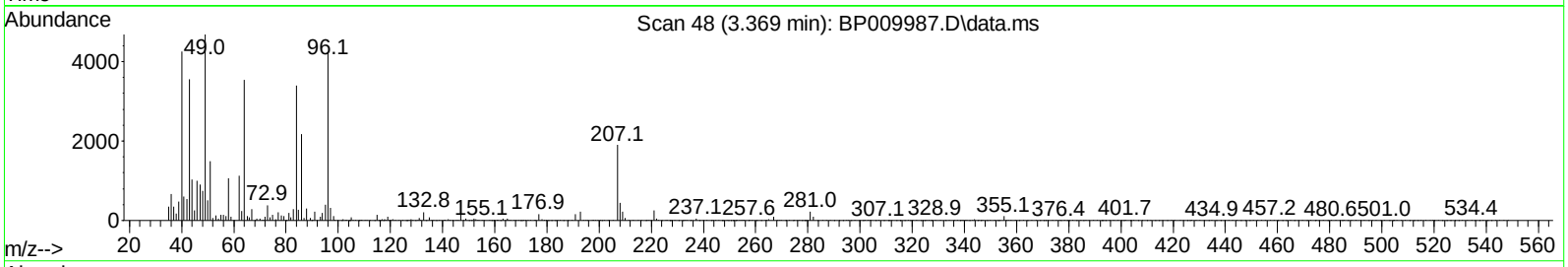

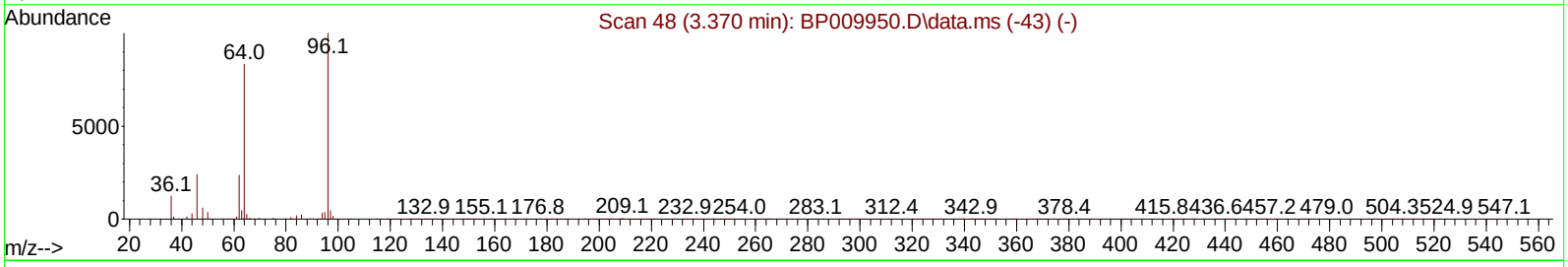

TIC: BP009987.D\data.ms

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

3.369min (-0.000) 2.20 ng/uL m

response 6111

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	84.04#
34.00	0.00	0.00
0.00	0.00	0.00

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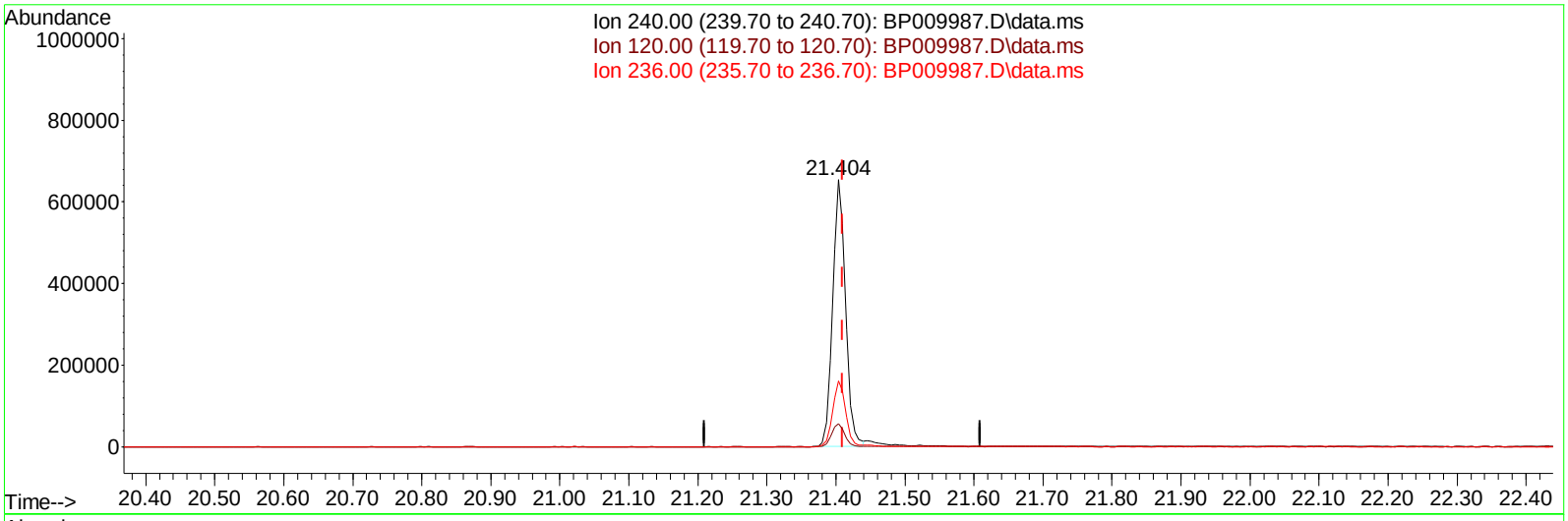

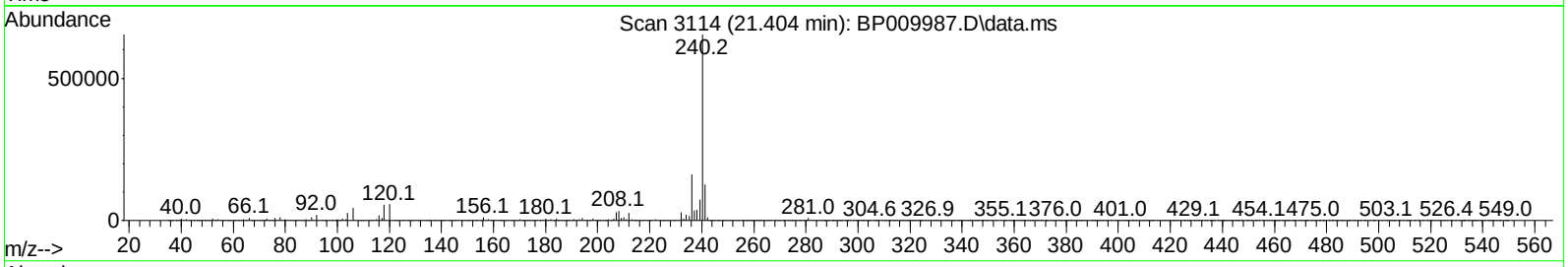

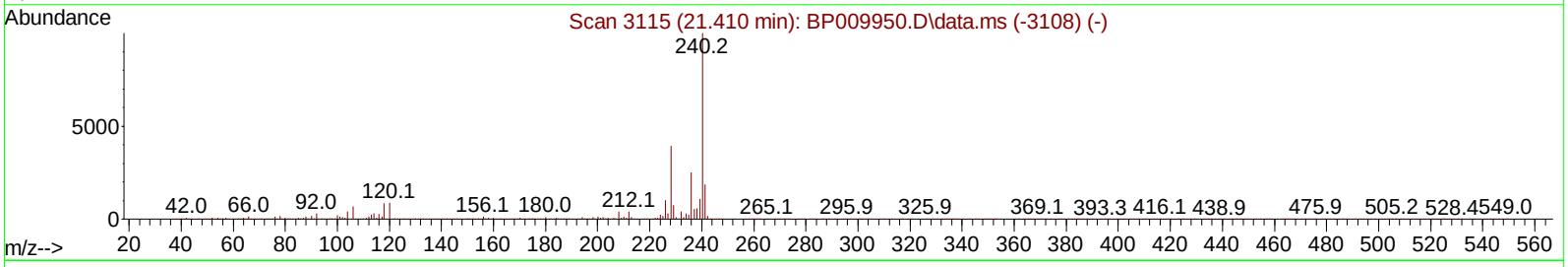

TIC: BP009987.D\data.ms

(79) Chrysene-d12 (I)

21.404min (-0.006) 20.00 ng/ul

response 853125

Ion	Exp%	Act%
240.00	100.00	100.00
120.00	9.30	8.68
236.00	31.60	24.61#
0.00	0.00	0.00

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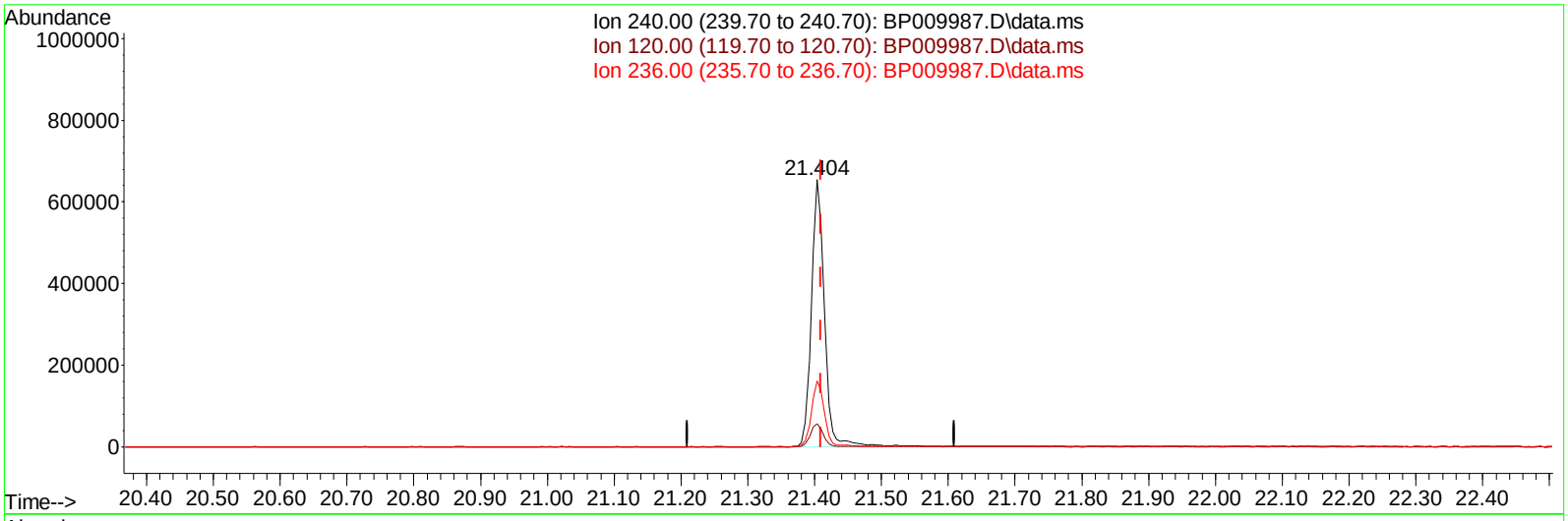

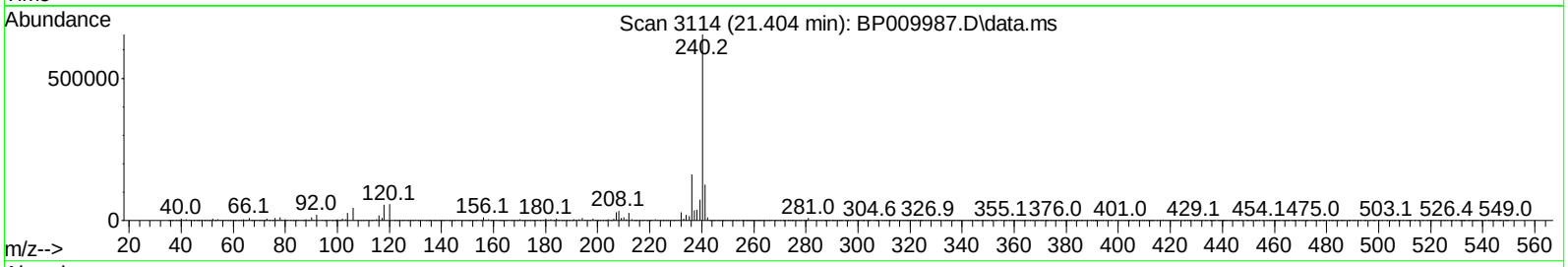

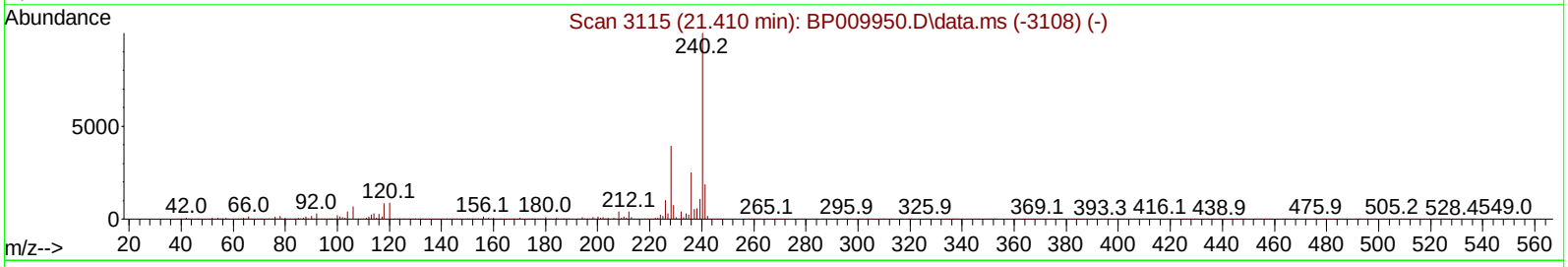

TIC: BP009987.D\data.ms

(79) Chrysene-d12 (I)

21.404min (-0.006) 20.00 ng/ul m

response 882388

Ion	Exp%	Act%
240.00	100.00	100.00
120.00	9.30	8.68
236.00	31.60	24.61#
0.00	0.00	0.00

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	123922	20.000	ng/ul	# 0.00
20) Naphthalene-d8	10.746	136	539925	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.575	164	389351	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.322	188	881462	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.404	240	882388m	20.000	ng/ul	0.00
88) Perylene-d12	23.857	264	746708	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	6111m	2.201	ng/uL	0.00
4) Pyridine-d5	3.787	84	74458	9.726	ng/ul	0.00
7) Phenol-d5	7.105	99	37987	3.492	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	207827	32.087	ng/ul	0.00
11) 2-Chlorophenol-d4	7.475	132	113710	13.856	ng/ul	0.00
15) 4-Methylphenol-d8	8.646	113	88800	9.777	ng/ul	0.00
21) Nitrobenzene-d5	9.099	128	139074	35.719	ng/ul	0.00
24) 2-Nitrophenol-d4	9.828	143	82078	19.228	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.369	165	142803	15.987	ng/ul	0.00
31) 4-Chloroaniline-d4	10.881	131	327976	25.779	ng/ul	0.00
46) Dimethylphthalate-d6	13.981	166	1079612	35.300	ng/ul	0.00
49) Acenaphthylene-d8	14.269	160	1174479	32.414	ng/ul	0.00
54) 4-Nitrophenol-d4	14.757	143	12062	2.175	ng/ul	0.00
60) Fluorene-d10	15.569	176	970941	37.738	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	68425	13.104	ng/ul	0.00
73) Anthracene-d10	17.422	188	1690647	40.059	ng/ul	0.00
81) Pyrene-d10	19.651	212	2079792	43.156	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.704	264	1563959	40.219	ng/ul	0.00

Target Compounds Qvalue

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

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