

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111523\
 Data File : BP018181.D
 Acq On : 15 Nov 2023 19:23
 Operator : MA/JU
 Sample : 05338-08
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCDL3

Quant Time: Nov 15 22:00:49 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 15 18:13:37 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	55603	20.000	ng/ul	0.00
20) Naphthalene-d8	10.598	136	222927	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.463	164	129919	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.245	188	274589	20.000	ng/ul	0.00
79) Chrysene-d12	21.380	240	252760	20.000	ng/ul	0.00
88) Perylene-d12	23.851	264	303505	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	4729	3.000	ng/uL	0.00
4) Pyridine-d5	3.587	84	96	0.023	ng/ul	0.00
7) Phenol-d5	6.963	99	27034	5.036	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.134	67	83920	26.602	ng/ul	0.00
11) 2-Chlorophenol-d4	7.310	132	82132	19.321	ng/ul	0.00
15) 4-Methylphenol-d8	8.516	113	53545	12.173	ng/ul	0.00
21) Nitrobenzene-d5	8.993	128	49244	24.518	ng/ul	0.00
24) 2-Nitrophenol-d4	9.710	143	53976	23.722	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.234	165	77710	19.162	ng/ul	0.00
31) 4-Chloroaniline-d4	10.793	131	75156	13.514	ng/ul	0.00
46) Dimethylphthalate-d6	13.887	166	324469	27.068	ng/ul	0.00
49) Acenaphthylene-d8	14.163	160	333516	25.544	ng/ul	0.00
54) 4-Nitrophenol-d4	14.728	143	16	0.009	ng/ul	0.00
60) Fluorene-d10	15.469	176	268885	27.169	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.634	200	33584	16.918	ng/ul	0.00
73) Anthracene-d10	17.345	188	448810	30.221	ng/ul	0.00
81) Pyrene-d10	19.604	212	520434	30.825	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.680	264	534624	30.294	ng/ul	0.00

Target Compounds Qvalue

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

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