

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111523\
 Data File : BP018183.D
 Acq On : 15 Nov 2023 20:31
 Operator : MA/JU
 Sample : 05338-10
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCDL5

Quant Time: Nov 15 22:01:11 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	57146	20.000	ng/ul	0.00
20) Naphthalene-d8	10.599	136	212597	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.463	164	123964	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.245	188	259892	20.000	ng/ul	0.00
79) Chrysene-d12	21.380	240	249961	20.000	ng/ul	0.00
88) Perylene-d12	23.851	264	302292	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	6088	3.758	ng/uL	0.00
4) Pyridine-d5	3.605	84	26929	6.375	ng/ul	0.01
7) Phenol-d5	6.963	99	28857	5.231	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.134	67	90431	27.892	ng/ul	0.00
11) 2-Chlorophenol-d4	7.305	132	90849	20.794	ng/ul	0.00
15) 4-Methylphenol-d8	8.516	113	55079	12.183	ng/ul	0.00
21) Nitrobenzene-d5	8.993	128	51598	26.939	ng/ul	0.00
24) 2-Nitrophenol-d4	9.705	143	55476	25.566	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.234	165	86915	22.473	ng/ul	0.00
31) 4-Chloroaniline-d4	10.793	131	73752	13.906	ng/ul	0.00
46) Dimethylphthalate-d6	13.887	166	344664	30.134	ng/ul	0.00
49) Acenaphthylene-d8	14.163	160	343684	27.588	ng/ul	0.00
54) 4-Nitrophenol-d4	14.740	143	12	0.007	ng/ul	0.00
60) Fluorene-d10	15.469	176	277739	29.411	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.781	200	122	0.065	ng/ul	0.15
73) Anthracene-d10	17.345	188	436316	31.041	ng/ul	0.00
81) Pyrene-d10	19.604	212	548794	32.869	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.680	264	589538	33.540	ng/ul	0.00

Target Compounds Qvalue

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

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