

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111523\
 Data File : BP018179.D
 Acq On : 15 Nov 2023 18:15
 Operator : MA/JU
 Sample : 05338-06
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCDL1

Quant Time: Nov 15 22:00:27 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	59870	20.000	ng/ul	0.00
20) Naphthalene-d8	10.599	136	224945	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.463	164	135349	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.245	188	306326	20.000	ng/ul	0.00
79) Chrysene-d12	21.380	240	304504	20.000	ng/ul	0.00
88) Perylene-d12	23.851	264	358532	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	5359	3.158	ng/uL	0.00
4) Pyridine-d5	3.581	84	70	0.016	ng/ul	-0.01
7) Phenol-d5	6.969	99	26582	4.599	ng/ul	0.02
9) Bis-(2-Chloroethyl)eth...	7.134	67	83961	24.718	ng/ul	0.00
11) 2-Chlorophenol-d4	7.310	132	83963	18.344	ng/ul	0.00
15) 4-Methylphenol-d8	8.516	113	50579	10.679	ng/ul	0.00
21) Nitrobenzene-d5	8.993	128	49778	24.562	ng/ul	0.00
24) 2-Nitrophenol-d4	9.710	143	52802	22.998	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.234	165	80129	19.581	ng/ul	0.00
31) 4-Chloroaniline-d4	10.799	131	42323	7.542	ng/ul	0.01
46) Dimethylphthalate-d6	13.887	166	356373	28.536	ng/ul	0.00
49) Acenaphthylene-d8	14.163	160	337883	24.841	ng/ul	0.00
54) 4-Nitrophenol-d4	14.734	143	8	0.004	ng/ul	0.00
60) Fluorene-d10	15.469	176	285170	27.658	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.722	200	500	0.226	ng/ul	0.09
73) Anthracene-d10	17.345	188	510033	30.785	ng/ul	0.00
81) Pyrene-d10	19.604	212	622251	30.593	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.686	264	652369	31.292	ng/ul	0.00

Target Compounds Qvalue

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

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