

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081822\
 Data File : BP011481.D
 Acq On : 18 Aug 2022 14:38
 Operator : CG/JU
 Sample : PB147077BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Aug 18 23:16:05 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Aug 17 01:58:37 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.564	152	106012	20.000	ng/u1	0.00
20) Naphthalene-d8	10.352	136	427082	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.240	164	227184	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.010	188	431999	20.000	ng/u1	0.00
79) Chrysene-d12	21.145	240	385406	20.000	ng/u1	0.01
88) Perylene-d12	23.445	264	412905	20.000	ng/u1	0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.070	96	19656	5.754	ng/uL	0.00
4) Pyridine-d5	3.481	84	240140	27.456	ng/u1	0.00
7) Phenol-d5	6.746	99	257498	26.432	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.911	67	167311	26.435	ng/u1	0.00
11) 2-Chlorophenol-d4	7.093	132	208714	29.129	ng/u1	0.00
15) 4-Methylphenol-d8	8.281	113	196829	25.242	ng/u1	0.00
21) Nitrobenzene-d5	8.728	128	95730	31.135	ng/u1	0.00
24) 2-Nitrophenol-d4	9.446	143	100890	32.242	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	9.981	165	159841	27.673	ng/u1	0.00
31) 4-Chloroaniline-d4	10.505	131	251574	26.164	ng/u1	0.00
46) Dimethylphthalate-d6	13.663	166	477896	29.741	ng/u1	0.00
49) Acenaphthylene-d8	13.928	160	534607	29.935	ng/u1	0.00
54) 4-Nitrophenol-d4	14.469	143	68875	22.247	ng/u1	0.00
60) Fluorene-d10	15.246	176	383100	28.340	ng/u1	0.01
65) 4,6-Dinitro-2-methylph...	15.381	200	60699	26.003	ng/u1	0.01
73) Anthracene-d10	17.110	188	579981	30.314	ng/u1	0.00
81) Pyrene-d10	19.375	212	613222	31.448	ng/u1	0.01
92) Benzo(a)pyrene-d12	23.298	264	615607	30.650	ng/u1	0.02

Target Compounds	Qvalue

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

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