

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110822\
 Data File : BP012580.D
 Acq On : 09 Nov 2022 09:05
 Operator : CG/JU
 Sample : PB148832BL
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Nov 09 10:13:06 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 08 22:40:23 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.769	152	228628	20.000	ng/ul	0.00
20) Naphthalene-d8	10.569	136	1172191	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.428	164	860956	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.186	188	1965127	20.000	ng/ul	0.00
79) Chrysene-d12	21.286	240	1800702	20.000	ng/ul	0.00
88) Perylene-d12	23.668	264	1647633	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.228	96	30725	5.411	ng/uL	0.00
4) Pyridine-d5	3.652	84	222626	13.649	ng/ul	0.00
7) Phenol-d5	6.963	99	649389	31.668	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.110	67	359272	29.349	ng/ul	0.00
11) 2-Chlorophenol-d4	7.310	132	484593	32.360	ng/ul	0.00
15) 4-Methylphenol-d8	8.499	113	547497	33.758	ng/ul	0.00
21) Nitrobenzene-d5	8.940	128	255752	33.645	ng/ul	0.00
24) 2-Nitrophenol-d4	9.663	143	287678	36.462	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.210	165	535725	31.332	ng/ul	0.00
31) 4-Chloroaniline-d4	10.728	131	565944	22.230	ng/ul	0.00
46) Dimethylphthalate-d6	13.845	166	1812374	30.751	ng/ul	0.00
49) Acenaphthylene-d8	14.122	160	2002762	29.736	ng/ul	0.00
54) 4-Nitrophenol-d4	14.687	143	274761	24.632	ng/ul	0.00
60) Fluorene-d10	15.428	176	1591742	31.544	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.563	200	274250	26.805	ng/ul	0.00
73) Anthracene-d10	17.292	188	2681611	31.670	ng/ul	0.00
81) Pyrene-d10	19.533	212	3057319	33.817	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.515	264	2607588	32.390	ng/ul	0.00

Target Compounds Qvalue

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

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