

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
 Data File : BP013237.D
 Acq On : 22 Dec 2022 09:57
 Operator : CG/JU
 Sample : PB149758TB 10X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Dec 22 21:44:35 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 21 23:38:37 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	471937	20.000	ng/ul	0.00
20) Naphthalene-d8	10.763	136	1872288	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.587	164	1035262	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.345	188	2067031	20.000	ng/ul	0.00
79) Chrysene-d12	21.433	240	2109320	20.000	ng/ul	0.00
88) Perylene-d12	23.910	264	2428867	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.334	96	59038	5.349	ng/uL	0.00
4) Pyridine-d5	3.764	84	595626	18.996	ng/ul	0.00
7) Phenol-d5	7.105	99	948267	24.668	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	620003	26.737	ng/ul	0.00
11) 2-Chlorophenol-d4	7.475	132	743757	24.578	ng/ul	0.00
15) 4-Methylphenol-d8	8.657	113	713726	22.660	ng/ul	0.00
21) Nitrobenzene-d5	9.116	128	339457	24.677	ng/ul	0.00
24) 2-Nitrophenol-d4	9.840	143	370120	23.899	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.381	165	697741	23.195	ng/ul	0.00
31) 4-Chloroaniline-d4	10.898	131	805065	18.958	ng/ul	0.00
46) Dimethylphthalate-d6	13.992	166	1892353	23.784	ng/ul	0.00
49) Acenaphthylene-d8	14.281	160	2170684	24.828	ng/ul	0.00
54) 4-Nitrophenol-d4	14.787	143	240198	16.783	ng/ul	0.00
60) Fluorene-d10	15.581	176	1686034	25.048	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.704	200	190534	14.324	ng/ul	0.00
73) Anthracene-d10	17.445	188	2528431	27.114	ng/ul	0.00
81) Pyrene-d10	19.674	212	3054544	25.229	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.745	264	3263700	26.958	ng/ul	0.00

Target Compounds Qvalue

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

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