

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
 Data File : BP013253.D
 Acq On : 22 Dec 2022 20:08
 Operator : CG/JU
 Sample : PB149781BL
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Dec 22 21:49:04 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	715220	20.000	ng/ul	0.00
20) Naphthalene-d8	10.763	136	3085714	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.587	164	1988680	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.345	188	4149857	20.000	ng/ul	0.00
79) Chrysene-d12	21.433	240	2661501	20.000	ng/ul	0.00
88) Perylene-d12	23.910	264	2907032	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.334	96	76245	4.558	ng/uL	0.00
4) Pyridine-d5	3.764	84	1157940	24.368	ng/ul	0.00
7) Phenol-d5	7.105	99	1480875	25.420	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	914023	26.009	ng/ul	0.00
11) 2-Chlorophenol-d4	7.475	132	1186337	25.869	ng/ul	0.00
15) 4-Methylphenol-d8	8.658	113	1208714	25.322	ng/ul	0.00
21) Nitrobenzene-d5	9.116	128	597550	26.357	ng/ul	0.00
24) 2-Nitrophenol-d4	9.840	143	686541	26.898	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.381	165	1267690	25.570	ng/ul	0.00
31) 4-Chloroaniline-d4	10.899	131	1693013	24.190	ng/ul	0.00
46) Dimethylphthalate-d6	13.998	166	3815606	24.965	ng/ul	0.00
49) Acenaphthylene-d8	14.281	160	4320152	25.724	ng/ul	0.00
54) 4-Nitrophenol-d4	14.787	143	512479	18.641	ng/ul	0.00
60) Fluorene-d10	15.581	176	3303126	25.546	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.704	200	462867	17.332	ng/ul	0.00
73) Anthracene-d10	17.445	188	4939062	26.381	ng/ul	0.00
81) Pyrene-d10	19.675	212	4939084	32.330	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.751	264	3842809	26.521	ng/ul	0.00

Target Compounds Qvalue

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

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