

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021125\
 Data File : BP024035.D
 Acq On : 11 Feb 2025 12:17
 Operator : RC/JU
 Sample : PB166618TB 10X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SLEB618

Quant Time: Feb 13 13:52:31 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Feb 13 12:56:03 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.763	152	400358	20.000	ng/u1	0.00
20) Naphthalene-d8	10.534	136	1678207	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.375	164	1065668	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.169	188	2219453	20.000	ng/u1	0.00
79) Chrysene-d12	21.598	240	2141264	20.000	ng/u1	0.00
88) Perylene-d12	24.951	264	2209299	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.287	96	63223	6.561	ng/uL	0.00
4) Pyridine-d5	3.699	84	915071	32.911	ng/u1	0.00
7) Phenol-d5	6.952	99	1185309	33.455	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.099	67	602114	32.106	ng/u1	0.00
11) 2-Chlorophenol-d4	7.305	132	975492	35.061	ng/u1	0.00
15) 4-Methylphenol-d8	8.469	113	949774	32.815	ng/u1	0.00
21) Nitrobenzene-d5	8.905	128	444226	33.085	ng/u1	0.00
24) 2-Nitrophenol-d4	9.622	143	474220	32.732	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.169	165	890067	33.489	ng/u1	0.01
31) 4-Chloroaniline-d4	10.669	131	1283310	32.287	ng/u1	0.00
46) Dimethylphthalate-d6	13.781	166	2554189	32.133	ng/u1	0.00
49) Acenaphthylene-d8	14.063	160	3024562	33.737	ng/u1	0.00
54) 4-Nitrophenol-d4	14.592	143	410126	25.980	ng/u1	0.00
60) Fluorene-d10	15.381	176	2318484	34.636	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.498	200	264520	19.319	ng/u1	0.00
73) Anthracene-d10	17.269	188	3746671	34.381	ng/u1	0.00
81) Pyrene-d10	19.633	212	4217081	36.756	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.721	264	4067744	35.288	ng/u1	0.00

Target Compounds	Qvalue

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

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