

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100124\
 Data File : BP022143.D
 Acq On : 01 Oct 2024 20:36
 Operator : RC/JU
 Sample : PB163717BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK717

Quant Time: Oct 02 07:08:25 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 24 15:23:13 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.699	152	162313	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.457	136	598400	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.310	164	447200	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.110	188	1086037	20.000	ng/u1	0.00
79) Chrysene-d12	21.539	240	1209221	20.000	ng/u1	0.00
88) Perylene-d12	24.821	264	1386748	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.217	96	18280	4.411	ng/uL	0.00
4) Pyridine-d5	3.634	84	257364	21.728	ng/u1	0.00
7) Phenol-d5	6.887	99	337600	25.427	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.034	67	197046	24.101	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.240	132	261340	27.205	ng/u1	0.00
15) 4-Methylphenol-d8	8.405	113	272137	25.866	ng/u1	0.00
21) Nitrobenzene-d5	8.834	128	122147	27.272	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.552	143	149429	29.425	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.087	165	293066	26.574	ng/u1	0.00
31) 4-Chloroaniline-d4	10.593	131	338090	25.693	ng/u1	-0.01
46) Dimethylphthalate-d6	13.710	166	906106	26.211	ng/u1	-0.01
49) Acenaphthylene-d8	13.998	160	977842	26.717	ng/u1	0.00
54) 4-Nitrophenol-d4	14.528	143	145178	24.552	ng/u1	0.00
60) Fluorene-d10	15.322	176	793316	25.944	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.434	200	147868	21.930	ng/u1	-0.01
73) Anthracene-d10	17.204	188	1342454	26.478	ng/u1	-0.01
81) Pyrene-d10	19.598	212	1755502	28.022	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.598	264	1956974	26.831	ng/u1	-0.01

Target Compounds Qvalue

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

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