

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011624\
 Data File : BP019385.D
 Acq On : 16 Jan 2024 10:42
 Operator : MA/JU
 Sample : PB158454BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK454

Quant Time: Jan 17 00:34:57 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jan 09 03:05:24 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.769	152	164374	20.000	ng/u1	-0.02
20) Naphthalene-d8	10.569	136	625648	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.416	164	472224	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.175	188	1180222	20.000	ng/u1	-0.01
79) Chrysene-d12	21.280	240	1170661	20.000	ng/u1	0.00
88) Perylene-d12	23.639	264	1163882	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.217	96	22235	5.115	ng/uL	0.00
4) Pyridine-d5	3.634	84	318175	25.312	ng/u1	0.00
7) Phenol-d5	6.964	99	389333	24.782	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.116	67	245506	26.126	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.311	132	284522	26.364	ng/u1	0.00
15) 4-Methylphenol-d8	8.499	113	318159	25.527	ng/u1	-0.02
21) Nitrobenzene-d5	8.940	128	152599	28.992	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.658	143	182168	29.786	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.199	165	321764	26.042	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.722	131	399052	25.568	ng/u1	0.00
46) Dimethylphthalate-d6	13.840	166	1127587	27.028	ng/u1	0.00
49) Acenaphthylene-d8	14.110	160	1190647	26.624	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.657	143	149032	23.263	ng/u1	-0.01
60) Fluorene-d10	15.416	176	971430	28.441	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.557	200	198232	22.750	ng/u1	0.00
73) Anthracene-d10	17.275	188	1598095	27.149	ng/u1	-0.01
81) Pyrene-d10	19.522	212	2048783	26.839	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.486	264	1795650	28.374	ng/u1	0.00

Target Compounds Qvalue

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

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