

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011224\
 Data File : BP019323.D
 Acq On : 12 Jan 2024 09:48
 Operator : MA/JU
 Sample : PB158377BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK377

Quant Time: Jan 12 18:11:20 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	7.775	152	104247	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.575	136	425309	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.428	164	318271	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.186	188	816753	20.000	ng/u1	0.00
79) Chrysene-d12	21.292	240	810584	20.000	ng/u1	0.00
88) Perylene-d12	23.645	264	797463	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.216	96	18975	6.883	ng/uL	0.00
4) Pyridine-d5	3.634	84	279234	35.027	ng/u1	0.00
7) Phenol-d5	6.963	99	283454	28.449	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.116	67	182035	30.545	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.310	132	204056	29.814	ng/u1	0.00
15) 4-Methylphenol-d8	8.504	113	223448	28.268	ng/u1	-0.01
21) Nitrobenzene-d5	8.945	128	107136	29.942	ng/u1	0.00
24) 2-Nitrophenol-d4	9.663	143	127168	30.588	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.204	165	235753	28.068	ng/u1	0.00
31) 4-Chloroaniline-d4	10.728	131	313732	29.570	ng/u1	0.00
46) Dimethylphthalate-d6	13.845	166	878600	31.247	ng/u1	0.00
49) Acenaphthylene-d8	14.122	160	902880	29.955	ng/u1	0.00
54) 4-Nitrophenol-d4	14.663	143	125263	29.011	ng/u1	0.00
60) Fluorene-d10	15.422	176	747526	32.472	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.563	200	156395	25.936	ng/u1	0.00
73) Anthracene-d10	17.286	188	1263442	31.016	ng/u1	0.00
81) Pyrene-d10	19.533	212	1621328	30.675	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.498	264	1396316	32.201	ng/u1	0.01

Target Compounds Qvalue

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

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