

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
 Data File : BP012477.D
 Acq On : 03 Nov 2022 06:53
 Operator : CG/JU
 Sample : PB148677BL
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK677

Quant Time: Nov 03 07:26:10 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 03 00:09:04 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	337028	20.000	ng/u1	0.00
20) Naphthalene-d8	10.581	136	1538129	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.434	164	1070257	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.198	188	2366398	20.000	ng/u1	0.00
79) Chrysene-d12	21.292	240	2110815	20.000	ng/u1	0.00
88) Perylene-d12	23.674	264	1844481	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	39856	4.762	ng/uL	0.00
4) Pyridine-d5	3.658	84	314653	13.086	ng/u1	0.00
7) Phenol-d5	6.958	99	766280	25.349	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.122	67	449238	24.895	ng/u1	0.00
11) 2-Chlorophenol-d4	7.316	132	588393	26.654	ng/u1	0.00
15) 4-Methylphenol-d8	8.499	113	634481	26.538	ng/u1	0.00
21) Nitrobenzene-d5	8.952	128	305823	30.661	ng/u1	0.00
24) 2-Nitrophenol-d4	9.669	143	352925	34.090	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.210	165	635140	28.309	ng/u1	0.00
31) 4-Chloroaniline-d4	10.734	131	1060383	31.741	ng/u1	0.00
46) Dimethylphthalate-d6	13.851	166	2052125	28.010	ng/u1	0.00
49) Acenaphthylene-d8	14.128	160	2229323	26.626	ng/u1	0.00
54) 4-Nitrophenol-d4	14.663	143	336135	24.241	ng/u1	0.00
60) Fluorene-d10	15.434	176	1769180	28.204	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.569	200	203309	16.502	ng/u1	0.00
73) Anthracene-d10	17.292	188	2949844	28.931	ng/u1	-0.01
81) Pyrene-d10	19.539	212	3344243	31.556	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.521	264	2687271	29.818	ng/u1	0.00

Target Compounds Qvalue

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

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