

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120123\
 Data File : BP018537.D
 Acq On : 03 Dec 2023 03:56
 Operator : MA/JU
 Sample : PB157193BL
 Misc :
 ALS Vial : 68 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK193

Quant Time: Dec 03 23:48:56 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 01 21:56:43 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.857	152	405471	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.657	136	1798858	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.487	164	1224310	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.233	188	2631748	20.000	ng/u1	0.00
79) Chrysene-d12	21.316	240	1786344	20.000	ng/u1	-0.01
88) Perylene-d12	23.692	264	1733290	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	74198	6.236	ng/uL	0.00
4) Pyridine-d5	3.687	84	1053572	33.057	ng/u1	0.00
7) Phenol-d5	7.028	99	1269845	31.207	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.193	67	795055	32.967	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.387	132	964128	30.366	ng/u1	-0.01
15) 4-Methylphenol-d8	8.569	113	1028647	31.237	ng/u1	-0.01
21) Nitrobenzene-d5	9.022	128	491655	30.708	ng/u1	0.00
24) 2-Nitrophenol-d4	9.740	143	512644	30.884	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.275	165	992521	29.242	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.793	131	1526905	34.423	ng/u1	-0.01
46) Dimethylphthalate-d6	13.904	166	3557171	32.554	ng/u1	0.00
49) Acenaphthylene-d8	14.181	160	3798480	31.569	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.687	143	543906	30.324	ng/u1	0.00
60) Fluorene-d10	15.481	176	2934559	31.994	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.592	200	401778	25.461	ng/u1	-0.01
73) Anthracene-d10	17.333	188	4479151	32.374	ng/u1	0.00
81) Pyrene-d10	19.569	212	4700657	33.805	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.533	264	3395043	33.593	ng/u1	-0.02

Target Compounds	Qvalue

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

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