

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120723\
 Data File : BP018694.D
 Acq On : 08 Dec 2023 03:38
 Operator : MA/JU
 Sample : PB157420BL
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK420

Quant Time: Dec 08 04:15:47 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Dec 04 21:39:40 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.846	152	155394	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.640	136	565347	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.475	164	347486	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.222	188	858364	20.000	ng/u1	0.00
79) Chrysene-d12	21.310	240	1088223	20.000	ng/u1	0.00
88) Perylene-d12	23.674	264	1561391	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.264	96	30973	6.792	ng/uL	0.00
4) Pyridine-d5	3.681	84	390890	32.002	ng/u1	0.00
7) Phenol-d5	7.011	99	404188	25.919	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.181	67	230301	24.917	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.375	132	309645	25.447	ng/u1	-0.01
15) 4-Methylphenol-d8	8.558	113	238402	18.890	ng/u1	-0.01
21) Nitrobenzene-d5	9.005	128	108988	21.660	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.728	143	111725	21.416	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.263	165	225028	21.095	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.781	131	321450	23.058	ng/u1	-0.01
46) Dimethylphthalate-d6	13.887	166	695660	22.431	ng/u1	-0.02
49) Acenaphthylene-d8	14.169	160	760128	22.258	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.675	143	123620	24.283	ng/u1	-0.01
60) Fluorene-d10	15.463	176	609248	23.403	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.581	200	90989	17.679	ng/u1	-0.01
73) Anthracene-d10	17.322	188	1078332	23.896	ng/u1	0.00
81) Pyrene-d10	19.557	212	1489769	17.587	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.522	264	2144032	23.550	ng/u1	-0.01

Target Compounds Qvalue

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

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