

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
 Data File : BP018119.D
 Acq On : 12 Nov 2023 23:58
 Operator : MA/JU
 Sample : PB156922BL
 Misc :
 ALS Vial : 100 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK922

Quant Time: Nov 13 03:24:41 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 10 21:55:18 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.840	152	101156	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.663	136	406889	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.516	164	233567	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.298	188	469262	20.000	ng/u1	-0.01
79) Chrysene-d12	21.427	240	428482	20.000	ng/u1	0.00
88) Perylene-d12	23.927	264	495799	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.222	96	29089	10.144	ng/uL	-0.01
4) Pyridine-d5	3.658	84	402511	53.835	ng/u1	0.00
7) Phenol-d5	7.010	99	498238	51.018	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.193	67	316582	55.162	ng/u1	0.00
11) 2-Chlorophenol-d4	7.363	132	404898	52.356	ng/u1	-0.01
15) 4-Methylphenol-d8	8.563	113	411400	51.409	ng/u1	-0.01
21) Nitrobenzene-d5	9.051	128	200937	54.813	ng/u1	0.00
24) 2-Nitrophenol-d4	9.757	143	234296	56.415	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.287	165	379241	51.235	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.839	131	520108	51.238	ng/u1	-0.01
46) Dimethylphthalate-d6	13.939	166	1174975	54.521	ng/u1	-0.01
49) Acenaphthylene-d8	14.216	160	1320975	56.278	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.775	143	144584	44.898	ng/u1	0.00
60) Fluorene-d10	15.522	176	966915	54.344	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.675	200	167598	49.404	ng/u1	-0.01
73) Anthracene-d10	17.398	188	1452446	57.229	ng/u1	-0.01
81) Pyrene-d10	19.651	212	1652715	57.745	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.762	264	1724516	59.818	ng/u1	-0.01

Target Compounds Qvalue

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

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