

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
 Data File : BP018096.D
 Acq On : 12 Nov 2023 09:51
 Operator : MA/JU
 Sample : PB156817BL
 Misc :
 ALS Vial : 76 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK817

Quant Time: Nov 12 15:39:28 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	7.846	152	101371	20.000	ng/u1	0.00
20) Naphthalene-d8	10.663	136	441271	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.522	164	296882	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.298	188	627319	20.000	ng/u1	-0.01
79) Chrysene-d12	21.428	240	415857	20.000	ng/u1	0.00
88) Perylene-d12	23.927	264	442960	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.228	96	16035	5.580	ng/uL	0.00
4) Pyridine-d5	3.664	84	215798	28.801	ng/u1	0.00
7) Phenol-d5	7.017	99	288256	29.454	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.193	67	180490	31.382	ng/u1	0.00
11) 2-Chlorophenol-d4	7.364	132	230113	29.692	ng/u1	-0.01
15) 4-Methylphenol-d8	8.569	113	247337	30.842	ng/u1	0.00
21) Nitrobenzene-d5	9.052	128	118646	29.843	ng/u1	0.00
24) 2-Nitrophenol-d4	9.763	143	138723	30.800	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.293	165	235402	29.324	ng/u1	0.00
31) 4-Chloroaniline-d4	10.846	131	328421	29.833	ng/u1	0.00
46) Dimethylphthalate-d6	13.940	166	837274	30.566	ng/u1	-0.01
49) Acenaphthylene-d8	14.216	160	884573	29.648	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.798	143	65887	16.096	ng/u1	0.02
60) Fluorene-d10	15.522	176	670519	29.648	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	64697	14.266	ng/u1	0.00
73) Anthracene-d10	17.398	188	1009547	29.756	ng/u1	-0.01
81) Pyrene-d10	19.657	212	1027541	36.992	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.763	264	789069	30.635	ng/u1	-0.01

Target Compounds Qvalue

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

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