

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
 Data File : BP018097.D
 Acq On : 12 Nov 2023 10:25
 Operator : MA/JU
 Sample : PB156819BL
 Misc :
 ALS Vial : 77 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK819

Quant Time: Nov 12 15:40:26 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	89747	20.000	ng/u1	0.00
20) Naphthalene-d8	10.663	136	377326	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.522	164	229671	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.298	188	427443	20.000	ng/u1	-0.01
79) Chrysene-d12	21.427	240	355450	20.000	ng/u1	0.00
88) Perylene-d12	23.927	264	421124	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.228	96	14837	5.832	ng/uL	0.00
4) Pyridine-d5	3.658	84	201114	30.318	ng/u1	0.00
7) Phenol-d5	7.010	99	259483	29.948	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.193	67	160063	31.435	ng/u1	0.00
11) 2-Chlorophenol-d4	7.369	132	208035	30.320	ng/u1	0.00
15) 4-Methylphenol-d8	8.569	113	210572	29.658	ng/u1	0.00
21) Nitrobenzene-d5	9.052	128	103240	30.369	ng/u1	0.00
24) 2-Nitrophenol-d4	9.763	143	117012	30.382	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.293	165	189546	27.614	ng/u1	0.00
31) 4-Chloroaniline-d4	10.846	131	266084	28.267	ng/u1	0.00
46) Dimethylphthalate-d6	13.940	166	625777	29.530	ng/u1	-0.01
49) Acenaphthylene-d8	14.216	160	686929	29.762	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.804	143	49890	15.755	ng/u1	0.02
60) Fluorene-d10	15.522	176	496710	28.390	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	55657	18.011	ng/u1	0.00
73) Anthracene-d10	17.404	188	709910	30.708	ng/u1	0.00
81) Pyrene-d10	19.651	212	760179	32.017	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.763	264	797787	32.580	ng/u1	-0.01

Target Compounds Qvalue

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

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