

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071723\
 Data File : BP016260.D
 Acq On : 17 Jul 2023 18:57
 Operator : MA/JU
 Sample : PB153975BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Jul 18 00:01:44 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jul 14 23:53:26 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.963	152	135199	20.000	ng/ul	0.00
20) Naphthalene-d8	10.787	136	663675	20.000	ng/ul	0.01
38) Acenaphthene-d10	14.616	164	434859	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.381	188	1018076	20.000	ng/ul	0.00
79) Chrysene-d12	21.492	240	975916	20.000	ng/ul	0.02
88) Perylene-d12	23.980	264	995953	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	25414	6.652	ng/uL	0.00
4) Pyridine-d5	3.734	84	318177	33.391	ng/ul	0.00
7) Phenol-d5	7.163	99	409283	33.515	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.293	67	275870	33.408	ng/ul	0.00
11) 2-Chlorophenol-d4	7.499	132	322412	36.573	ng/ul	0.00
15) 4-Methylphenol-d8	8.710	113	348159	35.391	ng/ul	0.00
21) Nitrobenzene-d5	9.169	128	144077	31.467	ng/ul	0.02
24) 2-Nitrophenol-d4	9.899	143	179139	34.328	ng/ul	0.02
28) 2,4-Dichlorophenol-d3	10.481	165	242142	26.795	ng/ul	0.04
31) 4-Chloroaniline-d4	10.987	131	404579	27.932	ng/ul	0.03
46) Dimethylphthalate-d6	14.040	166	1108174	33.407	ng/ul	0.01
49) Acenaphthylene-d8	14.310	160	1229740	33.828	ng/ul	0.00
54) 4-Nitrophenol-d4	14.981	143	36582	5.404	ng/ul	0.05
60) Fluorene-d10	15.622	176	983863	35.440	ng/ul	0.01
65) 4,6-Dinitro-2-methylph...	15.816	200	105240	17.907	ng/ul	0.02
73) Anthracene-d10	17.486	188	1575112	35.370	ng/ul	0.00
81) Pyrene-d10	19.722	212	1959988	38.704	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.804	264	1705255	35.654	ng/ul	0.00

Target Compounds	Qvalue

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

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