

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071023\
 Data File : BP016166.D
 Acq On : 10 Jul 2023 19:15
 Operator : MA/JU
 Sample : 03385-12
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 YBW42

Quant Time: Jul 10 23:22:09 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jul 10 23:19:51 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.016	152	403199	20.000	ng/ul	0.00
20) Naphthalene-d8	10.846	136	1833807	20.000	ng/ul	# 0.01
38) Acenaphthene-d10	14.669	164	1208484	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.439	188	2672980	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.539	240	2504730	20.000	ng/ul	# 0.00
88) Perylene-d12	24.068	264	2368737	20.000	ng/ul	-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.346	96	16598	1.481	ng/uL	0.00
4) Pyridine-d5	3.811	84	107386	3.801	ng/ul	0.02
7) Phenol-d5	7.252	99	228912	6.635	ng/ul	0.05
9) Bis-(2-Chloroethyl)eth...	7.358	67	152480	7.144	ng/ul	0.01
11) 2-Chlorophenol-d4	7.575	132	211187	8.110	ng/ul	0.03
15) 4-Methylphenol-d8	8.805	113	198391	7.280	ng/ul	0.05
21) Nitrobenzene-d5	9.252	128	100550	7.175	ng/ul	0.04
24) 2-Nitrophenol-d4	10.010	143	43059	2.632	ng/ul	0.07
28) 2,4-Dichlorophenol-d3	10.616	165	121226	4.159	ng/ul	0.11
31) 4-Chloroaniline-d4	11.110	131	190992	5.101	ng/ul	0.09
46) Dimethylphthalate-d6	14.104	166	858529	9.191	ng/ul	0.02
49) Acenaphthylene-d8	14.369	160	887937	8.456	ng/ul	0.00
54) 4-Nitrophenol-d4	14.987	143	106	0.009	ng/ul	0.00
60) Fluorene-d10	15.692	176	739815	9.619	ng/ul	0.02
65) 4,6-Dinitro-2-methylph...	15.945	200	12750	0.776	ng/ul	0.08
73) Anthracene-d10	17.557	188	1190105	9.747	ng/ul	0.02
81) Pyrene-d10	19.780	212	1542563	9.432	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.898	264	1221224	10.220	ng/ul	-0.02

Target Compounds Qvalue

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

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