

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070823\
 Data File : BP016090.D
 Acq On : 08 Jul 2023 13:34
 Operator : MA/JU
 Sample : 03332-16
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 YBTY9

Quant Time: Jul 08 22:16:37 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Jul 08 10:11:43 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.034	152	207143	20.000	ng/ul	0.00
20) Naphthalene-d8	10.863	136	924404	20.000	ng/ul	# 0.01
38) Acenaphthene-d10	14.681	164	603947	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.445	188	1307376	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.545	240	1265373	20.000	ng/ul	# 0.00
88) Perylene-d12	24.074	264	1312863	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	16037	2.785	ng/uL	0.00
4) Pyridine-d5	3.811	84	901	0.062	ng/ul	0.00
7) Phenol-d5	7.252	99	230006	12.977	ng/ul	0.04
9) Bis-(2-Chloroethyl)eth...	7.369	67	136858	12.481	ng/ul	0.01
11) 2-Chlorophenol-d4	7.581	132	196153	14.661	ng/ul	0.02
15) 4-Methylphenol-d8	8.816	113	178127	12.722	ng/ul	0.05
21) Nitrobenzene-d5	9.252	128	93904	13.292	ng/ul	0.04
24) 2-Nitrophenol-d4	10.004	143	106688	12.939	ng/ul	0.06
28) 2,4-Dichlorophenol-d3	10.604	165	141597	9.637	ng/ul	0.10
31) 4-Chloroaniline-d4	11.099	131	188361	9.980	ng/ul	0.08
46) Dimethylphthalate-d6	14.104	166	790266	16.929	ng/ul	0.02
49) Acenaphthylene-d8	14.381	160	869509	16.570	ng/ul	0.00
54) 4-Nitrophenol-d4	15.081	143	11402	1.851	ng/ul	0.12
60) Fluorene-d10	15.692	176	678964	17.664	ng/ul	0.02
65) 4,6-Dinitro-2-methylph...	15.916	200	18342	2.281	ng/ul	0.06
73) Anthracene-d10	17.557	188	1073451	17.975	ng/ul	0.01
81) Pyrene-d10	19.780	212	1406365	17.021	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.904	264	1214569	18.340	ng/ul	0.00

Target Compounds Qvalue

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

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