

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071023\
 Data File : BP016163.D
 Acq On : 10 Jul 2023 17:33
 Operator : MA/JU
 Sample : 03385-09
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EX7Y4

Quant Time: Jul 10 23:21:39 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	8.040	152	250617	20.000	ng/ul	0.02
20) Naphthalene-d8	10.863	136	1219103	20.000	ng/ul	# 0.03
38) Acenaphthene-d10	14.681	164	820673	20.000	ng/ul	0.01
64) Phenanthrene-d10	17.451	188	1790283	20.000	ng/ul	# 0.02
79) Chrysene-d12	21.545	240	1734999	20.000	ng/ul	# 0.00
88) Perylene-d12	24.074	264	1765496	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.346	96	7959	1.143	ng/uL	0.00
4) Pyridine-d5	3.822	84	64092	3.650	ng/ul	0.03
7) Phenol-d5	7.299	99	50242	2.343	ng/ul	0.09
9) Bis-(2-Chloroethyl)eth...	7.363	67	101	0.008	ng/ul	0.02
11) 2-Chlorophenol-d4	7.622	132	104656	6.466	ng/ul	0.08
15) 4-Methylphenol-d8	8.857	113	40978	2.419	ng/ul	0.10
21) Nitrobenzene-d5	9.304	128	27750	2.978	ng/ul	0.09
24) 2-Nitrophenol-d4	10.093	143	26505	2.437	ng/ul	0.15
28) 2,4-Dichlorophenol-d3	10.698	165	29887	1.542	ng/ul	0.19
31) 4-Chloroaniline-d4	11.151	131	46623	1.873	ng/ul	0.13
46) Dimethylphthalate-d6	14.128	166	447168	7.050	ng/ul	0.05
49) Acenaphthylene-d8	14.381	160	484483	6.794	ng/ul	0.02
54) 4-Nitrophenol-d4	14.992	143	22	0.003	ng/ul	0.00
60) Fluorene-d10	15.710	176	407635	7.805	ng/ul	0.04
65) 4,6-Dinitro-2-methylph...	15.875	200	11	0.001	ng/ul	0.01
73) Anthracene-d10	17.592	188	581546	7.111	ng/ul	0.05
81) Pyrene-d10	19.792	212	806105	7.115	ng/ul	0.01
92) Benzo(a)pyrene-d12	23.910	264	632157	7.098	ng/ul	-0.01

Target Compounds Qvalue

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

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