

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
 Data File : BG058998.D
 Acq On : 12 Sep 2023 1:43
 Operator : MA/JU
 Sample : 04234-11
 Misc :
 ALS Vial : 98 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 YC1G6

Quant Time: Sep 12 02:17:18 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Sep 11 00:07:05 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.320	152	155465	20.000	ng/u1	0.00
20) Naphthalene-d8	11.170	136	719952	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.948	164	534080	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.698	188	1438493	20.000	ng/u1	# 0.00
79) Chrysene-d12	22.016	240	1359236	20.000	ng/u1	0.00
88) Perylene-d12	25.588	264	1536779	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.644	96	4385	1.199	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/u1	
7) Phenol-d5	7.433	99	81439	5.195	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.639	67	46172	5.023	ng/u1	0.00
11) 2-Chlorophenol-d4	7.839	132	51061	5.078	ng/u1	0.00
15) 4-Methylphenol-d8	8.996	113	43919	3.626	ng/u1	0.00
21) Nitrobenzene-d5	9.513	128	27842	5.331	ng/u1	0.00
24) 2-Nitrophenol-d4	10.236	143	29970	5.230	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.759	165	60767	4.990	ng/u1	0.00
31) 4-Chloroaniline-d4	11.305	131	38807	2.370	ng/u1	0.00
46) Dimethylphthalate-d6	14.343	166	230307	5.559	ng/u1	0.00
49) Acenaphthylene-d8	14.648	160	244863	5.308	ng/u1	0.00
54) 4-Nitrophenol-d4	15.107	143	30165	4.384	ng/u1	0.00
60) Fluorene-d10	15.935	176	213630	5.811	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	16.029	200	27401	3.881	ng/u1	0.00
73) Anthracene-d10	17.792	188	341353	5.345	ng/u1	0.00
81) Pyrene-d10	20.065	212	445185	6.183	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.324	264	423067	5.616	ng/u1	-0.02
Target Compounds						
16) Acetophenone	9.166	105	42384	2.111	ng/u1#	Qvalue 86

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

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