

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111124\
 Data File : BG063291.D
 Acq On : 11 Nov 2024 16:13
 Operator : RC/JU
 Sample : P4644-06
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0CQ8

Quant Time: Nov 11 16:16:41 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110624.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 06 15:45:52 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.835	152	101085	20.000	ng/u1	0.00
20) Naphthalene-d8	10.626	136	462746	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.475	164	411445	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.224	188	996193	20.000	ng/u1	0.00
79) Chrysene-d12	21.478	240	1013250	20.000	ng/u1	# 0.00
88) Perylene-d12	24.516	264	1076134	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	12578	5.626	ng/uL	0.00
4) Pyridine-d5	3.705	84	64714	8.758	ng/u1	0.00
7) Phenol-d5	7.013	99	104020	11.404	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.166	67	166653	31.149	ng/u1	0.00
11) 2-Chlorophenol-d4	7.371	132	189596	32.024	ng/u1	0.00
15) 4-Methylphenol-d8	8.540	113	184339	23.843	ng/u1	0.00
21) Nitrobenzene-d5	8.987	128	112976	34.727	ng/u1	0.00
24) 2-Nitrophenol-d4	9.710	143	139535	33.186	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.250	165	275422	33.379	ng/u1	0.00
31) 4-Chloroaniline-d4	10.761	131	193623	18.016	ng/u1	0.00
46) Dimethylphthalate-d6	13.881	166	1003743	31.144	ng/u1	0.00
49) Acenaphthylene-d8	14.163	160	1051135	33.322	ng/u1	0.00
54) 4-Nitrophenol-d4	14.686	143	46163	10.462	ng/u1	0.00
60) Fluorene-d10	15.468	176	922481	33.915	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.597	200	198702	31.737	ng/u1	0.00
73) Anthracene-d10	17.324	188	1651921	38.583	ng/u1	0.00
81) Pyrene-d10	19.622	212	2043422	40.223	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.310	264	2058990	39.629	ng/u1	0.00

Target Compounds Qvalue

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

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