

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110321\
 Data File : BG050858.D
 Acq On : 3 Nov 2021 20:21
 Operator : CG/JU
 Sample : M4412-08
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BG364

Quant Time: Nov 07 07:46:51 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 02 14:49:05 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.229	152	32003	20.000	ng/u1	0.00
20) Naphthalene-d8	11.055	136	140715	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.851	164	99687	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.595	188	237035	20.000	ng/u1	-0.01
79) Chrysene-d12	21.896	240	210131	20.000	ng/u1	-0.01
88) Perylene-d12	25.292	264	203733	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.588	96	1685	1.699	ng/uL	0.00
4) Pyridine-d5	4.022	84	9954	3.356	ng/u1	0.00
7) Phenol-d5	7.371	99	17525	5.133	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.542	67	51399	23.305	ng/u1	0.00
11) 2-Chlorophenol-d4	7.753	132	43917	18.561	ng/u1	-0.01
15) 4-Methylphenol-d8	8.928	113	32354	12.037	ng/u1	0.00
21) Nitrobenzene-d5	9.404	128	30071	25.147	ng/u1	0.00
24) 2-Nitrophenol-d4	10.127	143	32089	24.134	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.668	165	51517	23.001	ng/u1	0.00
31) 4-Chloroaniline-d4	11.196	131	1867	0.550	ng/u1	0.00
46) Dimethylphthalate-d6	14.246	166	250003	32.780	ng/u1	0.00
49) Acenaphthylene-d8	14.551	160	254347	26.768	ng/u1	0.00
54) 4-Nitrophenol-d4	15.045	143	9284	6.714	ng/u1	0.00
60) Fluorene-d10	15.838	176	207858	30.766	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.955	200	41116	28.611	ng/u1	0.00
73) Anthracene-d10	17.695	188	375323	33.491	ng/u1	-0.01
81) Pyrene-d10	19.974	212	460796	33.952	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.057	264	372890	33.108	ng/u1	-0.02

Target Compounds Qvalue

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

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