

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\
 Data File : BG050905.D
 Acq On : 9 Nov 2021 8:38
 Operator : CG/JU
 Sample : M4445-13
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BG387

Quant Time: Nov 09 08:40:14 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	36082	20.000	ng/u1	0.00
20) Naphthalene-d8	11.050	136	156763	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.851	164	102654	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.595	188	231331	20.000	ng/u1	-0.01
79) Chrysene-d12	21.890	240	206710	20.000	ng/u1	-0.02
88) Perylene-d12	25.286	264	202686	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.588	96	5325	4.763	ng/uL	0.00
4) Pyridine-d5	4.017	84	25987	7.770	ng/u1	0.00
7) Phenol-d5	7.372	99	22279	5.788	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.542	67	66053	26.564	ng/u1	0.00
11) 2-Chlorophenol-d4	7.754	132	54943	20.596	ng/u1	-0.01
15) 4-Methylphenol-d8	8.923	113	37878	12.499	ng/u1	-0.01
21) Nitrobenzene-d5	9.399	128	37507	28.155	ng/u1	-0.01
24) 2-Nitrophenol-d4	10.127	143	36193	24.434	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.668	165	58709	23.529	ng/u1	0.00
31) 4-Chloroaniline-d4	11.185	131	81456	21.557	ng/u1	0.00
46) Dimethylphthalate-d6	14.246	166	240508	30.624	ng/u1	0.00
49) Acenaphthylene-d8	14.546	160	290622	29.701	ng/u1	-0.01
54) 4-Nitrophenol-d4	15.045	143	6394	4.490	ng/u1	0.00
60) Fluorene-d10	15.838	176	210911	30.315	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.956	200	32668	23.293	ng/u1	0.00
73) Anthracene-d10	17.695	188	356480	32.594	ng/u1	-0.01
81) Pyrene-d10	19.969	212	422637	31.656	ng/u1	-0.01
92) Benzo(a)pyrene-d12	25.057	264	353178	31.520	ng/u1	-0.02

Target Compounds	Qvalue

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

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