

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\
 Data File : BG050884.D
 Acq On : 8 Nov 2021 17:34
 Operator : CG/JU
 Sample : M4492-06
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BG1S3

Quant Time: Nov 08 17:35:51 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	35279	20.000	ng/u1	0.00
20) Naphthalene-d8	11.055	136	156055	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.851	164	105747	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.594	188	238627	20.000	ng/u1	-0.01
79) Chrysene-d12	21.889	240	206476	20.000	ng/u1	-0.02
88) Perylene-d12	25.291	264	202969	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.587	96	6274	5.740	ng/uL	0.00
4) Pyridine-d5	4.016	84	26186	8.008	ng/u1	0.00
7) Phenol-d5	7.371	99	27377	7.274	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.542	67	72044	29.633	ng/u1	0.00
11) 2-Chlorophenol-d4	7.759	132	60526	23.205	ng/u1	0.00
15) 4-Methylphenol-d8	8.928	113	45313	15.293	ng/u1	0.00
21) Nitrobenzene-d5	9.398	128	40576	30.596	ng/u1	-0.01
24) 2-Nitrophenol-d4	10.127	143	43256	29.334	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.667	165	64375	25.916	ng/u1	0.00
31) 4-Chloroaniline-d4	11.184	131	98449	26.172	ng/u1	0.00
46) Dimethylphthalate-d6	14.245	166	262666	32.467	ng/u1	0.00
49) Acenaphthylene-d8	14.551	160	318819	31.630	ng/u1	0.00
54) 4-Nitrophenol-d4	15.044	143	8605	5.866	ng/u1	0.00
60) Fluorene-d10	15.838	176	231043	32.238	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.955	200	38861	26.861	ng/u1	0.00
73) Anthracene-d10	17.694	188	422527	37.451	ng/u1	-0.01
81) Pyrene-d10	19.968	212	480660	36.042	ng/u1	-0.01
92) Benzo(a)pyrene-d12	25.056	264	399474	35.602	ng/u1	-0.02

Target Compounds Qvalue

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

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