

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\
 Data File : BG050892.D
 Acq On : 8 Nov 2021 23:46
 Operator : CG/JU
 Sample : PB140481BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK481

Quant Time: Nov 09 01:23:00 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	37155	20.000	ng/u1	0.00
20) Naphthalene-d8	11.055	136	162704	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.851	164	109404	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.594	188	239941	20.000	ng/u1	-0.01
79) Chrysene-d12	21.889	240	211849	20.000	ng/u1	-0.02
88) Perylene-d12	25.291	264	208222	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.593	96	6791	5.899	ng/uL	0.00
4) Pyridine-d5	4.010	84	102316	29.709	ng/u1	0.00
7) Phenol-d5	7.371	99	111308	28.081	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.541	67	75655	29.547	ng/u1	0.00
11) 2-Chlorophenol-d4	7.759	132	79238	28.845	ng/u1	0.00
15) 4-Methylphenol-d8	8.922	113	84892	27.205	ng/u1	-0.01
21) Nitrobenzene-d5	9.398	128	42041	30.406	ng/u1	-0.01
24) 2-Nitrophenol-d4	10.127	143	47053	30.605	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.667	165	71557	27.631	ng/u1	0.00
31) 4-Chloroaniline-d4	11.184	131	114194	29.117	ng/u1	0.00
46) Dimethylphthalate-d6	14.245	166	263268	31.453	ng/u1	0.00
49) Acenaphthylene-d8	14.545	160	324517	31.119	ng/u1	-0.01
54) 4-Nitrophenol-d4	15.039	143	37853	24.942	ng/u1	0.00
60) Fluorene-d10	15.838	176	224722	30.307	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.955	200	33602	23.099	ng/u1	0.00
73) Anthracene-d10	17.694	188	367207	32.370	ng/u1	-0.01
81) Pyrene-d10	19.968	212	417622	30.521	ng/u1	-0.01
92) Benzo(a)pyrene-d12	25.056	264	350158	30.420	ng/u1	-0.02

Target Compounds Qvalue

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

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