

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
 Data File : BG050977.D
 Acq On : 11 Nov 2021 21:49
 Operator : CG/JU
 Sample : PB140655BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK655

Quant Time: Nov 12 02:28:45 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 11 12:40:48 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.229	152	30510	20.000	ng/ul	0.00
20) Naphthalene-d8	11.055	136	138287	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.851	164	94417	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.594	188	208437	20.000	ng/ul	0.00
79) Chrysene-d12	21.889	240	179415	20.000	ng/ul	0.00
88) Perylene-d12	25.285	264	173151	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.593	96	5504	5.822	ng/uL	0.00
4) Pyridine-d5	4.016	84	86234	30.493	ng/ul	0.00
7) Phenol-d5	7.371	99	101293	31.120	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.541	67	66613	31.682	ng/ul	0.00
11) 2-Chlorophenol-d4	7.753	132	72925	32.329	ng/ul	0.00
15) 4-Methylphenol-d8	8.928	113	80016	31.227	ng/ul	0.00
21) Nitrobenzene-d5	9.404	128	39723	33.802	ng/ul	0.00
24) 2-Nitrophenol-d4	10.127	143	45808	35.056	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.667	165	70247	31.914	ng/ul	0.00
31) 4-Chloroaniline-d4	11.184	131	108797	32.639	ng/ul	0.00
46) Dimethylphthalate-d6	14.245	166	257318	35.622	ng/ul	0.00
49) Acenaphthylene-d8	14.551	160	313482	34.833	ng/ul	0.00
54) 4-Nitrophenol-d4	15.039	143	42998	32.830	ng/ul	0.00
60) Fluorene-d10	15.838	176	219311	34.273	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.955	200	38093	30.144	ng/ul	0.00
73) Anthracene-d10	17.694	188	358472	36.376	ng/ul	0.00
81) Pyrene-d10	19.968	212	406258	35.058	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.050	264	333640	34.856	ng/ul	0.00

Target Compounds Qvalue

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

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