

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081922\
 Data File : BG054661.D
 Acq On : 19 Aug 2022 11:20
 Operator : CG/JU
 Sample : PB146668BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK668

Quant Time: Aug 20 02:45:36 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Aug 18 03:22:26 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.164	152	40877	20.000	ng/u1	0.00
20) Naphthalene-d8	10.990	136	171356	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.791	164	111817	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.541	188	250935	20.000	ng/u1	0.00
79) Chrysene-d12	21.836	240	240120	20.000	ng/u1	0.00
88) Perylene-d12	25.214	264	242390	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.510	96	5891	5.701	ng/uL	0.00
4) Pyridine-d5	3.933	84	70606	22.928	ng/u1	0.00
7) Phenol-d5	7.294	99	97945	28.182	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.476	67	65374	29.397	ng/u1	0.00
11) 2-Chlorophenol-d4	7.688	132	72892	29.655	ng/u1	0.00
15) 4-Methylphenol-d8	8.851	113	76818	29.260	ng/u1	0.00
21) Nitrobenzene-d5	9.333	128	35248	32.771	ng/u1	0.00
24) 2-Nitrophenol-d4	10.056	143	27529	28.587	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.590	165	70456	27.909	ng/u1	0.00
31) 4-Chloroaniline-d4	11.113	131	102890	28.390	ng/u1	0.00
46) Dimethylphthalate-d6	14.192	166	249355	32.455	ng/u1	0.00
49) Acenaphthylene-d8	14.492	160	302228	32.854	ng/u1	0.00
54) 4-Nitrophenol-d4	14.956	143	32443	26.720	ng/u1	0.00
60) Fluorene-d10	15.784	176	222431	31.927	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.884	200	20081	23.499	ng/u1	0.00
73) Anthracene-d10	17.635	188	360168	32.999	ng/u1	0.00
81) Pyrene-d10	19.920	212	422006	33.881	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.979	264	403748	33.919	ng/u1	0.00

Target Compounds Qvalue

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

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