

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081922\
 Data File : BG054663.D
 Acq On : 19 Aug 2022 13:36
 Operator : CG/JU
 Sample : PB146669BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK669

Quant Time: Aug 20 02:46:34 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	8.158	152	44437	20.000 ng/u1	0.00
20) Naphthalene-d8	10.984	136	182646	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.792	164	119039	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.541	188	266866	20.000 ng/u1	0.00
79) Chrysene-d12	21.836	240	256819	20.000 ng/u1	0.00
88) Perylene-d12	25.215	264	256765	20.000 ng/u1	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.511	96	6823	6.074 ng/uL	0.00
4) Pyridine-d5	3.928	84	76611	22.885 ng/u1	-0.01
7) Phenol-d5	7.295	99	106592	28.213 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.471	67	69934	28.929 ng/u1	-0.01
11) 2-Chlorophenol-d4	7.682	132	78308	29.306 ng/u1	0.00
15) 4-Methylphenol-d8	8.852	113	83580	29.285 ng/u1	0.00
21) Nitrobenzene-d5	9.333	128	37751	32.928 ng/u1	0.00
24) 2-Nitrophenol-d4	10.056	143	30173	29.396 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.591	165	76788	28.537 ng/u1	0.00
31) 4-Chloroaniline-d4	11.114	131	118943	30.791 ng/u1	0.00
46) Dimethylphthalate-d6	14.192	166	267105	32.656 ng/u1	0.00
49) Acenaphthylene-d8	14.492	160	324020	33.086 ng/u1	0.00
54) 4-Nitrophenol-d4	14.962	143	35089	27.146 ng/u1	0.00
60) Fluorene-d10	15.785	176	235850	31.799 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.885	200	21495	23.652 ng/u1	0.00
73) Anthracene-d10	17.636	188	383564	33.045 ng/u1	0.00
81) Pyrene-d10	19.921	212	455395	34.184 ng/u1	0.00
92) Benzo(a)pyrene-d12	24.980	264	427252	33.884 ng/u1	0.00

Target Compounds Qvalue

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

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