

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
 Data File : BG050228.D
 Acq On : 9 Sep 2021 11:53
 Operator : CG/JU
 Sample : M3608-09
 Misc :
 ALS Vial : 56 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DBKK5

Quant Time: Sep 09 13:36:33 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 07 18:32:38 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.943	152	57948	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.769	136	182480	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.611	164	122143	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.366	188	267012	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.636	240	283624	20.000	ng/ul	0.00
88) Perylene-d12	24.856	264	336099	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.308	96	2678	2.285	ng/uL	0.00
4) Pyridine-d5	3.755	84	19859	4.773	ng/ul	0.02
7) Phenol-d5	7.127	99	25893	5.755	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.268	67	54448	20.721	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.479	132	77557	20.770	ng/ul	0.00
15) 4-Methylphenol-d8	8.672	113	82469	23.312	ng/ul	0.00
21) Nitrobenzene-d5	9.118	128	57146	45.625	ng/ul	0.00
24) 2-Nitrophenol-d4	9.846	143	44075	28.939	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.399	165	72993	25.111	ng/ul	0.00
31) 4-Chloroaniline-d4	10.910	131	84295	23.537	ng/ul	0.00
46) Dimethylphthalate-d6	14.011	166	273516	29.657	ng/ul	0.00
49) Acenaphthylene-d8	14.305	160	315103	28.693	ng/ul	0.00
54) 4-Nitrophenol-d4	14.863	143	6897	6.198	ng/ul	0.01
60) Fluorene-d10	15.603	176	225913	29.501	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.744	200	41365	25.682	ng/ul	0.00
73) Anthracene-d10	17.466	188	393587	32.835	ng/ul	0.00
81) Pyrene-d10	19.757	212	466400	33.014	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.632	264	692477	38.921	ng/ul	0.00

Target Compounds Qvalue

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

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