

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
 Data File : BG050142.D
 Acq On : 3 Sep 2021 22:23
 Operator : CG/JU
 Sample : M3566-16
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DBKJ1

Manual Integrations
 APPROVED

mohammad
 9/7/2021 11:34:17 AM

Quant Time: Sep 04 01:40:19 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Aug 30 14:59:55 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.949	152	28567	20.000	ng/u1	0.00
20) Naphthalene-d8	10.769	136	118554	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.611	164	78126	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.366	188	167964	20.000	ng/u1	-0.01
79) Chrysene-d12	21.637	240	176851	20.000	ng/u1	-0.01
88) Perylene-d12	24.850	264	178248	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.309	96	1170m	2.112	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/u1	
7) Phenol-d5	7.127	99	23936	9.867	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.274	67	13224	9.812	ng/u1	0.00
11) 2-Chlorophenol-d4	7.485	132	18500m	10.192	ng/u1	0.00
15) 4-Methylphenol-d8	8.678	113	18517	9.693	ng/u1	0.00
21) Nitrobenzene-d5	9.124	128	9767	10.626	ng/u1	0.00
24) 2-Nitrophenol-d4	9.853	143	12352	11.286	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.411	165	18893	9.679	ng/u1	0.00
31) 4-Chloroaniline-d4	10.910	131	19956	7.511	ng/u1	-0.01
46) Dimethylphthalate-d6	14.012	166	65149	10.896	ng/u1	-0.01
49) Acenaphthylene-d8	14.305	160	79346	11.318	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.858	143	6490m	7.502	ng/u1	0.01
60) Fluorene-d10	15.604	176	57263	11.296	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.750	200	7307	6.762	ng/u1	0.00
73) Anthracene-d10	17.466	188	90501	11.999	ng/u1	-0.01
81) Pyrene-d10	19.757	212	116297	12.169	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.633	264	109700	11.600	ng/u1	-0.02

Target Compounds Qvalue

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

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