

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062224\
 Data File : BG061656.D
 Acq On : 22 Jun 2024 19:53
 Operator : MA/JU
 Sample : P2776-05
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A4B93

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 06/24/2024
 Supervised By :mohammad ahmed 06/24/2024

Quant Time: Jun 23 00:40:01 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jun 21 05:03:55 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.073	152	71466	20.000	ng/u1	0.00
20) Naphthalene-d8	10.893	136	345375	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.707	164	233038	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.445	188	504879	20.000	ng/u1	0.00
79) Chrysene-d12	21.710	240	423142	20.000	ng/u1	0.00
88) Perylene-d12	24.936	264	518766	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.473	96	8298	4.194	ng/uL	0.00
4) Pyridine-d5	3.890	84	116246	19.435	ng/u1	0.00
7) Phenol-d5	7.221	99	175795	22.659	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.403	67	109341	22.608	ng/u1	0.00
11) 2-Chlorophenol-d4	7.603	132	130544	24.520	ng/u1	0.00
15) 4-Methylphenol-d8	8.772	113	139466	23.253	ng/u1	0.00
21) Nitrobenzene-d5	9.242	128	65394	27.886	ng/u1	0.00
24) 2-Nitrophenol-d4	9.965	143	72686	30.603	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.500	165	144008	25.606	ng/u1	0.00
31) 4-Chloroaniline-d4	11.017	131	184539	21.673	ng/u1	0.00
46) Dimethylphthalate-d6	14.107	166	440659	24.572	ng/u1	0.00
49) Acenaphthylene-d8	14.401	160	508008	25.849	ng/u1	0.00
54) 4-Nitrophenol-d4	14.877	143	59814	19.421	ng/u1	0.00
60) Fluorene-d10	15.694	176	393334	25.054	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.799	200	39310	18.208	ng/u1	0.00
73) Anthracene-d10	17.544	188	590345	25.616	ng/u1	0.00
81) Pyrene-d10	19.824	212	638795	27.027	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.718	264	651932	26.312	ng/u1	0.00
Target Compounds					Qvalue	
72) Phenanthrene	17.486	178	30497	1.148	ng/u1#	96
80) Fluoranthene	19.489	202	84622	3.086	ng/u1#	94
82) Pyrene	19.848	202	72656	2.492	ng/u1	99
85) Benzo(a)anthracene	21.687	228	36606m	1.232	ng/u1	
87) Chrysene	21.751	228	39135	1.392	ng/u1	95
90) Benzo(b)fluoranthene	23.913	252	51994	1.667	ng/u1	97
93) Benzo(a)pyrene	24.789	252	36604m	1.219	ng/u1	

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

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