

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG050123\
 Data File : BG057272.D
 Acq On : 2 May 2023 10:17
 Operator : CG/JU
 Sample : 02419-23
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EW920

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 05/03/2023
 Supervised By : Jagrut Upadhyay 05/03/2023

Quant Time: May 03 01:24:41 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG042823.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon May 01 12:45:38 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.379	152	17999	20.000	ng/u1	0.00
20) Naphthalene-d8	11.216	136	73853	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.993	164	47983	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.737	188	97143	20.000	ng/u1	# 0.00
79) Chrysene-d12	22.054	240	85912	20.000	ng/u1	0.00
88) Perylene-d12	25.585	264	94065	20.000	ng/u1	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.668	96	811	1.983	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/u1	
7) Phenol-d5	7.521	99	21861	12.843	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.686	67	14590	14.573	ng/u1	0.00
11) 2-Chlorophenol-d4	7.909	132	17126	15.268	ng/u1	0.00
15) 4-Methylphenol-d8	9.084	113	3565	2.756	ng/u1	0.00
21) Nitrobenzene-d5	9.560	128	10030	17.054	ng/u1	0.00
24) 2-Nitrophenol-d4	10.288	143	11116	16.195	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.834	165	17449	13.177	ng/u1	0.00
31) 4-Chloroaniline-d4	11.351	131	13919	7.931	ng/u1	0.00
46) Dimethylphthalate-d6	14.377	166	68216	17.942	ng/u1	0.00
49) Acenaphthylene-d8	14.694	160	77782	18.202	ng/u1	0.00
54) 4-Nitrophenol-d4	15.193	143	1326m	2.396	ng/u1	0.02
60) Fluorene-d10	15.980	176	59786	16.880	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	0.000	200	0	0.000	ng/u1	
73) Anthracene-d10	17.831	188	80565	18.169	ng/u1	0.00
81) Pyrene-d10	20.104	212	97756	19.170	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.338	264	86469	18.043	ng/u1	0.02
Target Compounds						
					Qvalue	
72) Phenanthrene	17.778	178	11228	2.220	ng/u1#	86
80) Fluoranthene	19.769	202	24178	3.991	ng/u1#	97
82) Pyrene	20.134	202	24789	4.042	ng/u1#	97
85) Benzo(a)anthracene	22.037	228	13591	2.237	ng/u1#	85
87) Chrysene	22.107	228	14701	2.564	ng/u1#	77
90) Benzo(b)fluoranthene	24.451	252	15630m	2.399	ng/u1	
91) Benzo(k)fluoranthene	24.528	252	6671m	1.053	ng/u1	
93) Benzo(a)pyrene	25.420	252	11055m	1.966	ng/u1	
94) Indeno(1,2,3-cd)pyrene	29.650	276	8860m	1.260	ng/u1	
96) Benzo(g,h,i)perylene	30.937	276	7341m	1.290	ng/u1	

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

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