

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060724\
 Data File : BG061556.D
 Acq On : 8 Jun 2024 7:19
 Operator : MA/JU
 Sample : P2640-17
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A4C41

Manual Integrations APPROVED

Reviewed By :Yogesh Patel 06/10/2024
 Supervised By :mohammad ahmed 06/10/2024

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 ahmed
 06/10/2024

Quant Time: Jun 09 23:12:44 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jun 06 12:15:45 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.871	152	27841	20.000	ng/u1	0.00
20) Naphthalene-d8	10.662	136	105039	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.511	164	78338	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.260	188	185112	20.000	ng/u1	0.00
79) Chrysene-d12	21.520	240	200525	20.000	ng/u1	0.00
88) Perylene-d12	24.622	264	231628	20.000	ng/u1	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.330	96	2468	4.961	ng/uL	-0.03
4) Pyridine-d5	3.741	84	33385	18.863	ng/u1	-0.03
7) Phenol-d5	7.060	99	52481	22.954	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.196	67	29788	20.729	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.413	132	41490	24.520	ng/u1	0.00
15) 4-Methylphenol-d8	8.594	113	43648	22.389	ng/u1	0.00
21) Nitrobenzene-d5	9.029	128	21545	26.641	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.751	143	26288	27.774	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.304	165	52168	28.272	ng/u1	0.00
31) 4-Chloroaniline-d4	10.803	131	38538	15.964	ng/u1	0.00
46) Dimethylphthalate-d6	13.917	166	155829	25.648	ng/u1	0.00
49) Acenaphthylene-d8	14.199	160	170513	26.995	ng/u1	0.00
54) 4-Nitrophenol-d4	14.763	143	17348m	23.095	ng/u1	0.00
60) Fluorene-d10	15.504	176	138708	26.443	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.650	200	21138	18.880	ng/u1	0.00
73) Anthracene-d10	17.360	188	230209	28.208	ng/u1	0.00
81) Pyrene-d10	19.657	212	278767	27.075	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.411	264	311079	27.935	ng/u1	0.01
Target Compounds					Qvalue	
72) Phenanthrene	17.301	178	51750	5.760	ng/u1	98
74) Anthracene	17.395	178	9878m	1.063	ng/u1	
80) Fluoranthene	19.323	202	97800	7.918	ng/u1	98
82) Pyrene	19.687	202	91506	7.137	ng/u1	98
85) Benzo(a)anthracene	21.502	228	53642	3.877	ng/u1	93
87) Chrysene	21.561	228	52979	4.158	ng/u1	99
90) Benzo(b)fluoranthene	23.647	252	73759	5.348	ng/u1#	98
91) Benzo(k)fluoranthene	23.700	252	21697	1.550	ng/u1#	91
93) Benzo(a)pyrene	24.469	252	51721	3.950	ng/u1#	97
94) Indeno(1,2,3-cd)pyrene	28.130	276	34867m	2.202	ng/u1	
96) Benzo(g,h,i)perylene	29.234	276	34668m	2.752	ng/u1	

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

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