

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070323\
 Data File : BG057963.D
 Acq On : 3 Jul 2023 20:31
 Operator : CG/JU
 Sample : 03246-21
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DCGK8

Manual Integrations
 APPROVED

Reviewed By :Sohil Jodhani 07/04/2023
 Supervised By :mohammad ahmed 07/04/2023

Quant Time: Jul 04 04:05:00 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Jul 02 07:50:15 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.937	152	45041	20.000	ng/u1	0.00
20) Naphthalene-d8	10.740	136	220592	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.571	164	156710	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.309	188	337776	20.000	ng/u1	0.00
79) Chrysene-d12	21.563	240	266465	20.000	ng/u1	0.00
88) Perylene-d12	24.730	264	330308	20.000	ng/u1	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.378	96	5435	4.225	ng/uL	0.01
4) Pyridine-d5	3.784	84	16795	4.301	ng/u1	0.00
7) Phenol-d5	7.097	99	123964	26.950	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.262	67	79680	28.745	ng/u1	0.00
11) 2-Chlorophenol-d4	7.467	132	94117	29.667	ng/u1	0.00
15) 4-Methylphenol-d8	8.637	113	80119	23.084	ng/u1	0.00
21) Nitrobenzene-d5	9.095	128	50603	28.145	ng/u1	0.00
24) 2-Nitrophenol-d4	9.818	143	55892	29.579	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.358	165	102726	28.339	ng/u1	0.00
31) 4-Chloroaniline-d4	10.875	131	40926	7.447	ng/u1	0.00
46) Dimethylphthalate-d6	13.972	166	339637	27.468	ng/u1	0.00
49) Acenaphthylene-d8	14.259	160	402058	29.798	ng/u1	0.00
54) 4-Nitrophenol-d4	14.765	143	33996	15.562	ng/u1	0.00
60) Fluorene-d10	15.558	176	306676	29.205	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.675	200	41111	22.414	ng/u1	0.00
73) Anthracene-d10	17.409	188	449166	28.657	ng/u1	0.00
81) Pyrene-d10	19.700	212	507914	30.265	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.512	264	509571	30.280	ng/u1	0.01
Target Compounds					Qvalue	
72) Phenanthrene	17.356	178	210646	12.260	ng/u1	100
74) Anthracene	17.444	178	22404	1.289	ng/u1	98
80) Fluoranthene	19.365	202	575740	29.607	ng/u1	100
82) Pyrene	19.730	202	455312	23.260	ng/u1	99
85) Benzo(a)anthracene	21.545	228	250339	13.084	ng/u1	99
87) Chrysene	21.610	228	307431	17.141	ng/u1	96
90) Benzo(b)fluoranthene	23.725	252	465251	22.749	ng/u1	99
91) Benzo(k)fluoranthene	23.784	252	160488m	8.006	ng/u1	
93) Benzo(a)pyrene	24.577	252	289556	15.996	ng/u1	99
94) Indeno(1,2,3-cd)pyrene	28.325	276	241914	10.064	ng/u1	100
95) Dibenzo(a,h)anthracene	28.361	278	64507m	3.254	ng/u1	
96) Benzo(g,h,i)perylene	29.453	276	208859m	10.638	ng/u1	

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

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