

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071524\
 Data File : BG062071.D
 Acq On : 15 Jul 2024 16:38
 Operator : MA/JU
 Sample : P3100-23
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A4D51

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 07/16/2024
 Supervised By :mohammad ahmed 07/20/2024

Quant Time: Jul 16 02:52:44 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jul 11 12:11:32 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.042	152	43962	20.000	ng/u1	0.00
20) Naphthalene-d8	10.850	136	206420	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.669	164	145967	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.413	188	329973	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.679	240	305564	20.000	ng/u1	0.00
88) Perylene-d12	24.887	264	371120	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.435	96	2981m	2.548	ng/uL	0.00
4) Pyridine-d5	3.864	84	16565	4.605	ng/u1	0.00
7) Phenol-d5	7.202	99	89410	18.584	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.360	67	41537	13.634	ng/u1	0.00
11) 2-Chlorophenol-d4	7.572	132	60985	18.480	ng/u1	0.00
15) 4-Methylphenol-d8	8.747	113	73571	19.422	ng/u1	0.00
21) Nitrobenzene-d5	9.211	128	31315	19.947	ng/u1	0.00
24) 2-Nitrophenol-d4	9.928	143	38858	21.675	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.468	165	82579	23.881	ng/u1	0.00
31) 4-Chloroaniline-d4	10.985	131	3473m	0.692	ng/u1	0.00
46) Dimethylphthalate-d6	14.070	166	287119	23.926	ng/u1	0.00
49) Acenaphthylene-d8	14.364	160	302000	24.067	ng/u1	0.00
54) 4-Nitrophenol-d4	14.881	143	38792	20.242	ng/u1	0.00
60) Fluorene-d10	15.662	176	244978	24.249	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.780	200	36239	20.912	ng/u1	0.00
73) Anthracene-d10	17.513	188	407111	26.426	ng/u1	0.00
81) Pyrene-d10	19.799	212	372710	20.737	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.663	264	329003	17.875	ng/u1	0.00
Target Compounds					Qvalue	
52) Acenaphthene	14.728	153	17436	1.833	ng/u1	93
56) Dibenzofuran	15.063	168	29098	2.165	ng/u1	90
61) Fluorene	15.715	166	63900	5.590	ng/u1	92
72) Phenanthrene	17.460	178	807195	46.394	ng/u1	100
74) Anthracene	17.548	178	175920	9.784	ng/u1	99
77) Carbazole	17.819	167	34010	2.248	ng/u1	94
80) Fluoranthene	19.470	202	1356185m	63.760	ng/u1	
82) Pyrene	19.828	202	808226	36.786	ng/u1#	92
85) Benzo(a)anthracene	21.655	228	660172	29.840	ng/u1	99
87) Chrysene	21.720	228	569239	27.485	ng/u1	96
90) Benzo(b)fluoranthene	23.870	252	816006m	34.899	ng/u1	
91) Benzo(k)fluoranthene	23.929	252	233496	10.014	ng/u1	99
93) Benzo(a)pyrene	24.740	252	362567m	16.375	ng/u1	
94) Indeno(1,2,3-cd)pyrene	28.547	276	275924	9.752	ng/u1	96
95) Dibenzo(a,h)anthracene	28.594	278	116281m	4.976	ng/u1	
96) Benzo(g,h,i)perylene	29.693	276	179414m	7.970	ng/u1	

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

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