

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082422\
 Data File : BG054719.D
 Acq On : 24 Aug 2022 14:06
 Operator : CG/JU
 Sample : N4120-06DL 10X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BLDG-19CDL

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 08/25/2022
 Supervised By : Jagrut Upadhyay 08/25/2022

Quant Time: Aug 24 15:14:54 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 23 03:09:46 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.159	152	49491	20.000	ng	0.00
21) Naphthalene-d8	10.985	136	208475	20.000	ng	0.00
39) Acenaphthene-d10	14.793	164	132783	20.000	ng	0.00
64) Phenanthrene-d10	17.536	188	316005	20.000	ng	0.00
76) Chrysene-d12	21.831	240	307313	20.000	ng	0.00
86) Perylene-d12	25.204	264	324530	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.697	112	30858	11.749	ng	0.00
7) Phenol-d6	7.301	99	42852	11.645	ng	0.00
23) Nitrobenzene-d5	9.328	82	26444	7.229	ng	0.00
42) 2,4,6-Tribromophenol	16.273	330	20264	19.056	ng	0.00
45) 2-Fluorobiphenyl	13.412	172	64416	7.154	ng	0.00
79) Terphenyl-d14	20.133	244	118747	7.676	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	11.038	128	45337	4.041	ng	98
52) Acenaphthene	14.851	154	23142	3.030	ng	99
55) Dibenzofuran	15.186	168	35962	3.107	ng	95
58) Fluorene	15.833	166	39864	4.166	ng	97
71) Phenanthrene	17.578	178	489544	29.064	ng	100
72) Anthracene	17.672	178	140343	8.598	ng	99
73) Carbazole	17.936	167	62460	4.301	ng	98
75) Fluoranthene	19.581	202	655286	32.998	ng	100
78) Pyrene	19.945	202	563668	26.003	ng	99
80) Butylbenzylphthalate	20.821	149	15035	7.035	ng	93
81) Benzo(a)anthracene	21.808	228	310903	14.896	ng	98
83) Chrysene	21.878	228	252976	12.736	ng	98
84) Bis(2-ethylhexyl)phtha...	21.702	149	21643	6.026	ng	98
87) Indeno(1,2,3-cd)pyrene	29.088	276	170104	7.024	ng	# 95
88) Benzo(b)fluoranthene	24.129	252	323943	16.076	ng	99
89) Benzo(k)fluoranthene	24.193	252	119642m	5.753	ng	
90) Benzo(a)pyrene	25.045	252	248507	14.368	ng	97
91) Dibenzo(a,h)anthracene	29.152	278	43860m	2.205	ng	
92) Benzo(g,h,i)perylene	30.298	276	161289	8.293	ng	97

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

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