

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091922\
 Data File : BF130327.D
 Acq On : 19 Sep 2022 18:42
 Operator : CG\JU
 Sample : N4665-02DL 20X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMP-2DL

Quant Time: Sep 20 02:27:13 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 19 15:06:14 2022
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By : Christian Giraldo 09/20/2022
 Supervised By : Jagrut Upadhyay 09/20/2022

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.657	152	92752	20.000	ng	0.00
21) Naphthalene-d8	7.951	136	332269	20.000	ng	0.00
39) Acenaphthene-d10	9.692	164	165580	20.000	ng	0.00
64) Phenanthrene-d10	11.175	188	210422	20.000	ng	0.00
76) Chrysene-d12	13.810	240	196968	20.000	ng	0.00
86) Perylene-d12	15.198	264	222275	20.000	ng	# 0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.245	112	21502	4.021	ng	-0.01
7) Phenol-d6	6.287	99	27819	4.158	ng	-0.01
23) Nitrobenzene-d5	7.222	82	17659	2.715	ng	-0.01
42) 2,4,6-Tribromophenol	10.475	330	5095	3.211	ng	0.00
45) 2-Fluorobiphenyl	9.016	172	31029	2.729	ng	0.00
79) Terphenyl-d14	12.757	244	31737m	2.669	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	7.986	128	7256911	423.932	ng	94
37) 2-Methylnaphthalene	8.675	142	3672006	336.439	ng	99
38) 1-Methylnaphthalene	8.769	142	2469732	235.553	ng	99
46) 1,1'-Biphenyl	9.116	154	520432	42.085	ng	98
49) Acenaphthylene	9.557	152	644443	42.966	ng	# 91
52) Acenaphthene	9.722	154	244693	24.830	ng	97
55) Dibenzofuran	9.892	168	215005	15.686	ng	# 80
58) Fluorene	10.245	166	1144813	102.125	ng	99
71) Phenanthrene	11.216	178	4072004	344.677	ng	97
72) Anthracene	11.257	178	796131	69.832	ng	100
73) Carbazole	11.404	167	45356	4.408	ng	# 88
75) Fluoranthene	12.392	202	1432330	125.469	ng	96
78) Pyrene	12.621	202	2438358	141.511	ng	98
81) Benzo(a)anthracene	13.798	228	1148219m	87.628	ng	
83) Chrysene	13.839	228	1001713m	80.513	ng	
87) Indeno(1,2,3-cd)pyrene	16.521	276	284908	18.075	ng	96
88) Benzo(b)fluoranthene	14.815	252	789291m	54.405	ng	
89) Benzo(k)fluoranthene	14.833	252	214891m	15.058	ng	
90) Benzo(a)pyrene	15.145	252	923578	80.356	ng	98
91) Dibenzo(a,h)anthracene	16.527	278	103546	8.008	ng	# 90
92) Benzo(g,h,i)perylene	16.915	276	294622	22.618	ng	99

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

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