

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071119\
 Data File : BF115516.D
 Acq On : 12 Jul 2019 1:07
 Operator : HP/JU
 Sample : K3726-01DL 5X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PX-1_070919DL

Manual Integrations
 APPROVED

mohammad
 7/12/2019 10:37:54 PM

Quant Time: Jul 12 07:34:23 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF070519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 10 16:47:05 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	133537	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	522890	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	218650	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	362104	20.00	ng	0.00
76) Chrysene-d12	14.06	240	318817	20.00	ng	0.00
87) Perylene-d12	15.54	264	265635	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	174739	20.22	ng	-0.01
7) Phenol-d6	6.51	99	233413	21.23	ng	-0.01
23) Nitrobenzene-d5	7.45	82	127796	14.44	ng	-0.01
42) 2,4,6-Tribromophenol	10.72	330	43994	18.20	ng	-0.01
45) 2-Fluorobiphenyl	9.25	172	269671	21.64	ng	0.00
79) Terphenyl-d14	13.00	244	249588	16.03	ng	0.00
Target Compounds						
31) Naphthalene	8.20	128	55277	2.128	ng	98
37) 2-Methylnaphthalene	8.89	142	56804	3.272	ng	98
38) 1-Methylnaphthalene	8.99	142	70627	4.264	ng	97
49) Acenaphthylene	9.79	152	163347	7.594	ng	97
50) Dimethylphthalate	9.64	163	56457	3.440	ng	99
52) Acenaphthene	9.96	154	65309	4.824	ng	99
55) Dibenzofuran	10.14	168	84822	4.448	ng	# 94
58) Fluorene	10.48	166	207481	13.803	ng	99
71) Phenanthrene	11.45	178	1123701	59.006	ng	99
72) Anthracene	11.49	178	321440	16.832	ng	99
73) Carbazole	11.65	167	64811	3.710	ng	99
75) Fluoranthene	12.63	202	989948	49.370	ng	99
78) Pvrene	12.87	202	1216699	49.075	ng	100
81) Benzo(a)anthracene	14.05	228	584431	28.610	ng	95
83) Chrysene	14.09	228	506521	24.133	ng	96
86) Indeno(1,2,3-cd)pyrene	17.03	276	192442	6.101	ng	96
88) Benzo(b)fluoranthene	15.11	252	543053m	31.399	ng	
89) Benzo(k)fluoranthene	15.13	252	157994m	10.272	ng	
90) Benzo(a)pvrene	15.48	252	393628	26.387	ng	99
91) Dibenzo(a,h)anthracene	17.04	278	54009	4.237	ng	# 95
92) Benzo(g,h,i)perylene	17.47	276	173765	14.402	ng	98

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

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