

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072519\
 Data File : BF115755.D
 Acq On : 25 Jul 2019 12:38
 Operator : HP/JU
 Sample : K3887-03DL 20X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AOC-7(2)DL

Manual Integrations
 APPROVED

mohammad
 7/26/2019 3:02:42 PM

Quant Time: Jul 25 13:28:07 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 25 13:16:07 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	120055	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	509011	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	294929	20.00	ng	0.00
64) Phenanthrene-d10	11.39	188	532505	20.00	ng	0.00
76) Chrysene-d12	14.03	240	355912	20.00	ng	0.00
87) Perylene-d12	15.50	264	390277	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	26264	3.58	ng	0.00
7) Phenol-d6	6.50	99	49508	5.32	ng	0.00
23) Nitrobenzene-d5	7.43	82	31363	3.58	ng	0.00
42) 2,4,6-Tribromophenol	10.69	330	14528	4.22	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	62948	3.74	ng	-0.01
79) Terphenyl-d14	12.97	244	72891	3.78	ng	-0.01
Target Compounds						Qvalue
31) Naphthalene	8.17	128	1127240	45.727	ng	99
37) 2-Methylnaphthalene	8.86	142	587513	35.954	ng	99
38) 1-Methylnaphthalene	8.96	142	515673	33.224	ng	99
46) 1,1'-Biphenyl	9.33	154	127970	5.550	ng	99
49) Acenaphthylene	9.77	152	204567	7.191	ng	98
52) Acenaphthene	9.94	154	69703	3.795	ng	91
55) Dibenzofuran	10.11	168	55333	2.121	ng	# 91
58) Fluorene	10.45	166	465140	24.946	ng	100
71) Phenanthrene	11.42	178	1930387	70.721	ng	99
72) Anthracene	11.47	178	295773	10.485	ng	99
75) Fluoranthene	12.60	202	980356	33.988	ng	97
78) Pylene	12.84	202	1492140	49.484	ng	100
81) Benzo(a)anthracene	14.02	228	545347	23.258	ng	93
83) Chrysene	14.06	228	564614	23.987	ng	99
86) Indeno(1,2,3-cd)pyrene	16.97	276	205139	10.258	ng	97
88) Benzo(b)fluoranthene	15.07	252	384458m	15.298	ng	
89) Benzo(k)fluoranthene	15.09	252	129177m	5.765	ng	
90) Benzo(a)pyrene	15.44	252	397219	18.390	ng	98
91) Dibenzo(a,h)anthracene	16.97	278	65253	3.441	ng	94
92) Benzo(a,h,i)perylene	17.42	276	236770	12.520	ng	97

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

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