

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111919\
 Data File : BF117880.D
 Acq On : 20 Nov 2019 3:54
 Operator : JU
 Sample : K5908-03
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B-8-2-(6-12)

Manual Integrations
 APPROVED

mohammad
 11/21/2019 9:03:55 AM

Quant Time: Nov 20 05:43:41 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 13 18:36:07 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.92	152	156063	20.00	ng	-0.01
21) Naphthalene-d8	8.20	136	592886	20.00	ng	-0.01
39) Acenaphthene-d10	9.96	164	277832	20.00	ng	-0.01
64) Phenanthrene-d10	11.46	188	448824	20.00	ng	-0.01
76) Chrysene-d12	14.10	240	393289	20.00	ng	-0.01
87) Perylene-d12	15.62	264	313204	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	959017	108.77	ng	0.00
7) Phenol-d6	6.54	99	1225180	115.21	ng	0.00
23) Nitrobenzene-d5	7.47	82	751593	75.36	ng	-0.02
42) 2,4,6-Tribromophenol	10.76	330	284843	96.84	ng	-0.01
45) 2-Fluorobiphenyl	9.28	172	1399267	90.35	ng	-0.01
79) Terphenyl-d14	13.04	244	1474279	68.51	ng	-0.01
Target Compounds						
10) Phenol	6.55	94	67906	6.145	ng	83
50) Dimethylphthalate	9.67	163	256685	13.175	ng	99
58) Fluorene	10.52	166	39563	2.328	ng	97
71) Phenanthrene	11.48	178	547586	24.267	ng	97
72) Anthracene	11.53	178	101438	4.534	ng	99
73) Carbazole	11.69	167	54520	2.789	ng	99
75) Fluoranthene	12.67	202	661425	29.616	ng	99
78) Pyrene	12.90	202	488402	15.366	ng	99
81) Benzo(a)anthracene	14.09	228	263133	10.125	ng	99
83) Chrysene	14.13	228	236230	9.380	ng	97
86) Indeno(1,2,3-cd)pyrene	17.14	276	71122	3.318	ng	95
88) Benzo(b)fluoranthene	15.16	252	232306m	11.416	ng	
89) Benzo(k)fluoranthene	15.19	252	64174m	3.347	ng	
90) Benzo(a)pyrene	15.54	252	127822	7.143	ng	96
92) Benzo(a,h,i)perylene	17.60	276	58544	3.816	ng	97

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

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