

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061119\
 Data File : BF114960.D
 Acq On : 11 Jun 2019 22:54
 Operator : HP/JU
 Sample : K3247-03
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-GF-02-041-B

Manual Integrations
 APPROVED

mohammad
 6/12/2019 9:09:05 AM

Quant Time: Jun 12 05:30:03 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 10 16:44:00 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.66	152	274409	20.00	ng	0.00
21) Naphthalene-d8	7.94	136	1041373	20.00	ng	0.00
39) Acenaphthene-d10	9.69	164	477262	20.00	ng	0.00
64) Phenanthrene-d10	11.16	188	797603	20.00	ng	0.00
76) Chrysene-d12	13.79	240	601607	20.00	ng	0.00
87) Perylene-d12	15.17	264	504011	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.27	112	1572985	97.01	ng	0.00
7) Phenol-d6	6.31	99	1951114	99.76	ng	0.00
23) Nitrobenzene-d5	7.22	82	1186520	70.01	ng	0.00
42) 2,4,6-Tribromophenol	10.47	330	431645	83.14	ng	0.00
45) 2-Fluorobiphenyl	9.02	172	2004153	66.65	ng	0.00
79) Terphenyl-d14	12.74	244	1710586	56.55	ng	0.00
Target Compounds						
10) Phenol	6.32	94	54205	2.486	ng	Qvalue # 73
50) Dimethylphthalate	9.41	163	304711	8.495	ng	97
71) Phenanthrene	11.19	178	322143	8.325	ng	95
72) Anthracene	11.23	178	85468	2.176	ng	95
75) Fluoranthene	12.37	202	616631	15.406	ng	95
78) Pyrene	12.59	202	665088	12.694	ng	96
81) Benzo(a)anthracene	13.77	228	366274	8.234	ng	97
83) Chrysene	13.81	228	319751	7.158	ng	97
86) Indeno(1,2,3-cd)pyrene	16.47	276	106085	2.526	ng	93
88) Benzo(b)fluoranthene	14.78	252	323138m	9.848	ng	
89) Benzo(k)fluoranthene	14.80	252	102463m	3.598	ng	
90) Benzo(a)pyrene	15.11	252	209190	7.340	ng	97
92) Benzo(a,h,i)perylene	16.86	276	94898	3.955	ng	97

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

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