

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061119\
 Data File : BF114963.D
 Acq On : 12 Jun 2019 00:13
 Operator : HP/JU
 Sample : K3248-03
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-GF-02-042-B

Manual Integrations
 APPROVED

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

Quant Time: Jun 12 05:36:42 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	270125	20.00	ng	0.00
21) Naphthalene-d8	7.94	136	1000308	20.00	ng	0.00
39) Acenaphthene-d10	9.69	164	454179	20.00	ng	0.00
64) Phenanthrene-d10	11.16	188	786583	20.00	ng	0.00
76) Chrysene-d12	13.79	240	559571	20.00	ng	0.00
87) Perylene-d12	15.17	264	448429	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.27	112	1914036	119.92	ng	0.00
7) Phenol-d6	6.31	99	2370999	123.16	ng	0.00
23) Nitrobenzene-d5	7.23	82	1440767	88.50	ng	0.00
42) 2,4,6-Tribromophenol	10.47	330	555336	112.40	ng	0.00
45) 2-Fluorobiphenyl	9.02	172	2381621	83.22	ng	0.00
79) Terphenyl-d14	12.75	244	2259972	80.32	ng	0.00
Target Compounds						
10) Phenol	6.32	94	62825	2.927	ng	Qvalue # 66
50) Dimethylphthalate	9.41	163	363679	10.654	ng	98
71) Phenanthrene	11.19	178	331471	8.686	ng	96
72) Anthracene	11.23	178	90351	2.333	ng	97
75) Fluoranthene	12.37	202	581011	14.719	ng	94
78) Pyrene	12.59	202	535105	10.980	ng	96
81) Benzo(a)anthracene	13.77	228	286619	6.927	ng	96
83) Chrysene	13.81	228	253911	6.111	ng	96
86) Indeno(1,2,3-cd)pyrene	16.47	276	101729	2.604	ng	93
88) Benzo(b)fluoranthene	14.79	252	272304m	9.327	ng	
89) Benzo(k)fluoranthene	14.80	252	94739m	3.739	ng	
90) Benzo(a)pyrene	15.11	252	178048	7.022	ng	# 95
92) Benzo(a,h,i)perylene	16.87	276	99170	4.645	ng	# 91

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

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