

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061119\
 Data File : BF114951.D
 Acq On : 11 Jun 2019 18:56
 Operator : HP/JU
 Sample : K3163-07 2X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SU-04-060719-B

Manual Integrations
 APPROVED

mohammad
 6/12/2019 9:04:02 AM

Quant Time: Jun 12 04:52:21 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	248280	20.00	ng	0.00
21) Naphthalene-d8	7.94	136	945906	20.00	ng	0.00
39) Acenaphthene-d10	9.69	164	450578	20.00	ng	0.00
64) Phenanthrene-d10	11.16	188	789450	20.00	ng	0.00
76) Chrysene-d12	13.79	240	604174	20.00	ng	0.00
87) Perylene-d12	15.17	264	536999	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.26	112	657053	44.79	ng	0.00
7) Phenol-d6	6.30	99	823589	46.54	ng	-0.01
23) Nitrobenzene-d5	7.22	82	484547	31.47	ng	-0.01
42) 2,4,6-Tribromophenol	10.47	330	188594	38.48	ng	0.00
45) 2-Fluorobiphenyl	9.02	172	984976	34.69	ng	0.00
79) Terphenyl-d14	12.75	244	904275	29.77	ng	0.00
Target Compounds						Qvalue
49) Acenaphthylene	9.55	152	129256	3.082	ng	98
50) Dimethylphthalate	9.41	163	105952	3.129	ng	97
71) Phenanthrene	11.19	178	395507	10.326	ng	97
72) Anthracene	11.24	178	121830	3.134	ng	97
75) Fluoranthene	12.37	202	694682	17.535	ng	95
78) Pyrene	12.60	202	797020	15.147	ng	97
81) Benzo(a)anthracene	13.78	228	467573m	10.467	ng	
83) Chrysene	13.82	228	479588	10.691	ng	97
86) Indeno(1,2,3-cd)pyrene	16.49	276	202780	4.808	ng	95
88) Benzo(b)fluoranthene	14.79	252	557533m	15.948	ng	
89) Benzo(k)fluoranthene	14.81	252	173815m	5.729	ng	
90) Benzo(a)pyrene	15.12	252	363857	11.982	ng	96
91) Dibenzo(a,h)anthracene	16.49	278	53030	2.067	ng	# 76
92) Benzo(g,h,i)perylene	16.88	276	181776	7.110	ng	99

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

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