

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031819\
 Data File : BF113089.D
 Acq On : 18 Mar 2019 14:43
 Operator : JU/SJ
 Sample : K1901-07
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-6(0.5-1.5)

Manual Integrations
 APPROVED

Sohil
 3/19/2019 5:09:41 PM

Quant Time: Mar 19 05:18:50 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 13 03:42:55 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	190250	20.00	ng	0.00
21) Naphthalene-d8	8.01	136	675419	20.00	ng	0.00
39) Acenaphthene-d10	9.76	164	324204	20.00	ng	-0.01
64) Phenanthrene-d10	11.24	188	603289	20.00	ng	0.00
76) Chrysene-d12	13.87	240	539266	20.00	ng	0.00
87) Perylene-d12	15.27	264	476322	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.35	112	953460	77.96	ng	0.00
7) Phenol-d6	6.37	99	1247485	81.20	ng	0.00
23) Nitrobenzene-d5	7.29	82	772128	68.41	ng	-0.01
42) 2,4,6-Tribromophenol	10.55	330	310927	90.34	ng	0.00
45) 2-Fluorobiphenyl	9.09	172	1166175	54.14	ng	0.00
79) Terphenyl-d14	12.82	244	1246848	44.82	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
10) Phenol	6.39	94	40188	2.242	ng	98
50) Dimethylphthalate	9.47	163	106886	4.471	ng	99
71) Phenanthrene	11.26	178	106120	3.535	ng	97
75) Fluoranthene	12.44	202	365255	11.358	ng	95
78) Pyrene	12.67	202	415315	9.136	ng	99
81) Benzo(a)anthracene	13.86	228	280756	7.512	ng	97
83) Chrysene	13.89	228	276983	7.566	ng	98
86) Indeno(1,2,3-cd)pyrene	16.63	276	140163m	4.491	ng	
88) Benzo(b)fluoranthene	14.87	252	396368m	12.865	ng	
89) Benzo(k)fluoranthene	14.90	252	121709m	4.361	ng	
90) Benzo(a)pyrene	15.22	252	219559	8.057	ng	98
92) Benzo(a,h,i)perylene	17.04	276	144130	6.172	ng	96

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

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