

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031219\
 Data File : BF113005.D
 Acq On : 12 Mar 2019 16:18
 Operator : JU/SJ
 Sample : K1687-03 2X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HR-06-031119-B

Manual Integrations
 APPROVED

Sohil
 3/13/2019 11:40:35 AM

Quant Time: Mar 13 04:19:56 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	184392	20.00	ng	0.00
21) Naphthalene-d8	8.01	136	643105	20.00	ng	0.00
39) Acenaphthene-d10	9.76	164	268733	20.00	ng	0.00
64) Phenanthrene-d10	11.24	188	457885	20.00	ng	0.00
76) Chrysene-d12	13.86	240	474778	20.00	ng	0.00
87) Perylene-d12	15.27	264	452245	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.34	112	460176	38.82	ng	0.00
7) Phenol-d6	6.36	99	597133	40.10	ng	-0.01
23) Nitrobenzene-d5	7.29	82	322394	30.00	ng	-0.01
42) 2,4,6-Tribromophenol	10.55	330	129089	45.25	ng	0.00
45) 2-Fluorobiphenyl	9.09	172	555664	28.41	ng	0.00
79) Terphenyl-d14	12.82	244	486134	19.85	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
50) Dimethylphthalate	9.48	163	48122	2.429	ng	# 98
71) Phenanthrene	11.26	178	226368	9.936	ng	98
72) Anthracene	11.32	178	84119	3.527	ng	99
75) Fluoranthene	12.45	202	327783	13.429	ng	97
78) Pyrene	12.67	202	313532	7.833	ng	99
81) Benzo(a)anthracene	13.86	228	166958	5.074	ng	95
83) Chrysene	13.89	228	160392	4.976	ng	97
86) Indeno(1,2,3-cd)pyrene	16.63	276	58515m	2.129	ng	
88) Benzo(b)fluoranthene	14.87	252	190084m	6.498	ng	
89) Benzo(k)fluoranthene	14.90	252	54770m	2.067	ng	
90) Benzo(a)pyrene	15.21	252	132713	5.129	ng	97
92) Benzo(a,h,i)perylene	17.04	276	52439	2.365	ng	96

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

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