

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030819\
 Data File : BF112951.D
 Acq On : 8 Mar 2019 18:22
 Operator : JU/SJ
 Sample : K1693-03 2X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OR-01-030719-B

Manual Integrations
 APPROVED

Sohil
 3/11/2019 9:32:22 AM

Quant Time: Mar 08 23:43:40 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 01 15:40:43 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	187645	20.00	ng	-0.01
21) Naphthalene-d8	8.01	136	703771	20.00	ng	-0.01
39) Acenaphthene-d10	9.76	164	306295	20.00	ng	-0.01
64) Phenanthrene-d10	11.23	188	517310	20.00	ng	-0.01
76) Chrysene-d12	13.86	240	483049	20.00	ng	-0.01
87) Perylene-d12	15.27	264	445317	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	556902	46.17	ng	-0.01
7) Phenol-d6	6.36	99	725556	47.88	ng	0.00
23) Nitrobenzene-d5	7.29	82	413726	35.18	ng	-0.02
42) 2,4,6-Tribromophenol	10.54	330	174599	53.69	ng	-0.02
45) 2-Fluorobiphenyl	9.09	172	762719	35.15	ng	-0.01
79) Terphenyl-d14	12.82	244	655884	26.32	ng	-0.01
Target Compounds						
37) 2-Methylnaphthalene	8.72	142	47606	2.110	ng	99
49) Acenaphthylene	9.62	152	69286	2.092	ng	99
50) Dimethylphthalate	9.48	163	53070	2.350	ng	# 95
71) Phenanthrene	11.26	178	172220	6.691	ng	97
75) Fluoranthene	12.44	202	245145	8.890	ng	98
78) Pyrene	12.67	202	357730	8.785	ng	99
81) Benzo(a)anthracene	13.85	228	144199m	4.307	ng	
83) Chrysene	13.89	228	167795	5.117	ng	97
88) Benzo(b)fluoranthene	14.87	252	155127m	5.386	ng	
90) Benzo(a)pyrene	15.20	252	119193	4.678	ng	94
92) Benzo(g,h,i)perylene	17.02	276	59603	2.730	ng	96

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

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