

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF041819\
 Data File : BF113835.D
 Acq On : 18 Apr 2019 23:54
 Operator : JU/SJ
 Sample : K2197-13
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-01-041619-E

Manual Integrations
 APPROVED

Sohil
 4/23/2019 8:16:51 AM

Quant Time: Apr 19 03:05:05 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF041819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 18 13:55:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	134810	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	444748	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	200479	20.00	ng	0.00
64) Phenanthrene-d10	11.41	188	431807	20.00	ng	0.00
76) Chrysene-d12	14.04	240	470868	20.00	ng	0.00
87) Perylene-d12	15.52	264	416288	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	693743	85.57	ng	0.00
7) Phenol-d6	6.52	99	838331	82.92	ng	0.00
23) Nitrobenzene-d5	7.45	82	471771	64.60	ng	0.00
42) 2,4,6-Tribromophenol	10.71	330	233727	105.52	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	854311	64.99	ng	0.00
79) Terphenyl-d14	12.99	244	1255810	55.93	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.64	163	88327	6.296	ng	99
71) Phenanthrene	11.43	178	51454	2.340	ng	97
75) Fluoranthene	12.62	202	137050	5.880	ng	95
78) Pyrene	12.85	202	129475	3.902	ng	96
81) Benzo(a)anthracene	14.03	228	77111	2.802	ng	93
83) Chrysene	14.07	228	76989	2.818	ng	96
88) Benzo(b)fluoranthene	15.08	252	101021m	3.893	ng	
90) Benzo(a)pyrene	15.45	252	57941	2.525	ng	# 89

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

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