

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF031919\
 Data File : BF113126.D
 Acq On : 19 Mar 2019 15:33
 Operator : JU/SJ
 Sample : K1970-04 20X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P-4-E.WALL-S.CENTER

Manual Integrations
 APPROVED

Sohil
 3/20/2019 6:51:21 PM

Quant Time: Mar 20 08:10:57 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	6.72	152	185962	20.00	ng	-0.01
21) Naphthalene-d8	8.00	136	561422	20.00	ng	-0.01
39) Acenaphthene-d10	9.76	164	267310	20.00	ng	0.00
64) Phenanthrene-d10	11.24	188	578443	20.00	ng	0.00
76) Chrysene-d12	13.86	240	416061	20.00	ng	0.00
87) Perylene-d12	15.27	264	424249	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	90238	7.55	ng	-0.02
7) Phenol-d6	6.36	99	148932	9.92	ng	-0.02
23) Nitrobenzene-d5	7.28	82	111896	11.93	ng	-0.02
42) 2,4,6-Tribromophenol	10.55	330	30165	10.63	ng	0.00
45) 2-Fluorobiphenyl	9.08	172	159025	6.64	ng	-0.01
79) Terphenyl-d14	12.82	244	160568	7.48	ng	0.00
Target Compounds						
37) 2-Methylnaphthalene	8.73	142	791667	43.981	ng	99
38) 1-Methylnaphthalene	8.82	142	592141	33.960	ng	100
52) Acenaphthene	9.79	154	48145	2.849	ng	# 81
55) Dibenzofuran	9.96	168	53734m	2.187	ng	
58) Fluorene	10.30	166	147299m	8.448	ng	
71) Phenanthrene	11.26	178	364162	12.653	ng	96

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

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