

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF042919\
 Data File : BF114027.D
 Acq On : 29 Apr 2019 17:18
 Operator : JU/SJ
 Sample : K2553-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RBR200034

Manual Integrations
 APPROVED

Sohil
 4/30/2019 6:42:20 PM

Quant Time: Apr 30 05:59:56 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF042219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 23 03:32:35 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	126126	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	471956	20.00	ng	0.00
39) Acenaphthene-d10	9.92	164	247724	20.00	ng	-0.01
64) Phenanthrene-d10	11.41	188	483222	20.00	ng	0.00
76) Chrysene-d12	14.04	240	319742	20.00	ng	-0.02
87) Perylene-d12	15.54	264	364220	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	664445	90.10	ng	0.00
7) Phenol-d6	6.52	99	869202	93.95	ng	0.00
23) Nitrobenzene-d5	7.45	82	525468	68.74	ng	0.00
42) 2,4,6-Tribromophenol	10.71	330	277976	92.11	ng	-0.01
45) 2-Fluorobiphenyl	9.25	172	1084724	68.48	ng	-0.01
79) Terphenyl-d14	12.99	244	1153874	73.98	ng	-0.01
Target Compounds						
50) Dimethylphthalate	9.63	163	162428	9.066	ng	99
75) Fluoranthene	12.62	202	101401	3.827	ng	97
78) Pyrene	12.85	202	85669	3.831	ng	96
83) Chrysene	14.07	228	42199	2.300	ng	97
88) Benzo(b)fluoranthene	15.10	252	58313m	2.531	ng	

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

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