

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040819\
 Data File : BF113649.D
 Acq On : 8 Apr 2019 19:58
 Operator : JU/SJ
 Sample : K2305-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-EF-02-001-ABCD

Manual Integrations
 APPROVED

Sohil
 4/9/2019 10:15:59 AM

Quant Time: Apr 09 01:47:50 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 04 04:05:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.68	152	120400	20.00	ng	0.00
21) Naphthalene-d8	7.96	136	460122	20.00	ng	0.00
39) Acenaphthene-d10	9.71	164	223319	20.00	ng	-0.01
64) Phenanthrene-d10	11.19	188	427217	20.00	ng	-0.01
76) Chrysene-d12	13.83	240	308613	20.00	ng	-0.02
87) Perylene-d12	15.23	264	295011	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	666102	94.34	ng	0.00
7) Phenol-d6	6.33	99	885901	101.80	ng	0.00
23) Nitrobenzene-d5	7.25	82	524330	66.75	ng	-0.01
42) 2,4,6-Tribromophenol	10.50	330	257101	95.84	ng	-0.01
45) 2-Fluorobiphenyl	9.04	172	977359	70.46	ng	-0.01
79) Terphenyl-d14	12.77	244	896286	59.13	ng	-0.02
Target Compounds						
50) Dimethylphthalate	9.43	163	91728	5.572	ng	99
71) Phenanthrene	11.22	178	44144	2.079	ng	98
75) Fluoranthene	12.40	202	78798	3.464	ng	100
78) Pyrene	12.63	202	70939	3.021	ng	98
88) Benzo(b)fluoranthene	14.84	252	43572m	2.413	ng	

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

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