

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF041619\
 Data File : BF113803.D
 Acq On : 16 Apr 2019 19:51
 Operator : JU/SJ
 Sample : K2411-03 2X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-16-S

Manual Integrations
 APPROVED

Sohil
 4/17/2019 3:45:44 PM

Quant Time: Apr 17 04:09:59 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	6.67	152	131756	20.00	ng	-0.02
21) Naphthalene-d8	7.96	136	490041	20.00	ng	-0.01
39) Acenaphthene-d10	9.70	164	200966	20.00	ng	-0.02
64) Phenanthrene-d10	11.19	188	354649	20.00	ng	-0.02
76) Chrysene-d12	13.83	240	346482	20.00	ng	-0.02
87) Perylene-d12	15.23	264	375331	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	373518	48.34	ng	0.00
7) Phenol-d6	6.33	99	507896	53.33	ng	0.00
23) Nitrobenzene-d5	7.24	82	280465	33.52	ng	-0.02
42) 2,4,6-Tribromophenol	10.50	330	99085	41.04	ng	-0.01
45) 2-Fluorobiphenyl	9.03	172	556398	44.57	ng	-0.02
79) Terphenyl-d14	12.76	244	505741	29.72	ng	-0.03
Target Compounds						
50) Dimethylphthalate	9.43	163	48784	3.293	ng	97
75) Fluoranthene	12.40	202	68952	3.651	ng	100
78) Pyrene	12.63	202	68201	2.587	ng	98
81) Benzo(a)anthracene	13.82	228	41914m	2.097	ng	
83) Chrysene	13.85	228	43474	2.146	ng	99
88) Benzo(b)fluoranthene	14.84	252	65693m	2.859	ng	
90) Benzo(a)pyrene	15.19	252	48207	2.361	ng	99

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

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