

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040219\
 Data File : BF113516.D
 Acq On : 2 Apr 2019 20:14
 Operator : JU/SJ
 Sample : K2180-11
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-6D

Manual Integrations
 APPROVED

Sohil
 4/3/2019 12:42:38 PM

Quant Time: Apr 03 07:17:29 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF032519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 26 03:32:08 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.69	152	138662	20.00	ng	-0.01
21) Naphthalene-d8	7.97	136	512572	20.00	ng	-0.01
39) Acenaphthene-d10	9.72	164	239081	20.00	ng	-0.02
64) Phenanthrene-d10	11.20	188	427573	20.00	ng	-0.02
76) Chrysene-d12	13.83	240	337929	20.00	ng	-0.02
87) Perylene-d12	15.23	264	374214	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.31	112	924377	109.00	ng	0.00
7) Phenol-d6	6.35	99	1190313	110.95	ng	-0.01
23) Nitrobenzene-d5	7.26	82	715773	82.88	ng	-0.01
42) 2,4,6-Tribromophenol	10.51	330	327778	110.99	ng	-0.02
45) 2-Fluorobiphenyl	9.05	172	1221161	78.76	ng	-0.02
79) Terphenyl-d14	12.78	244	1080992	70.51	ng	-0.02
Target Compounds						
50) Dimethylphthalate	9.44	163	86342	4.873	ng	99
71) Phenanthrene	11.22	178	118701	5.590	ng	99
75) Fluoranthene	12.41	202	117093	4.878	ng	98
78) Pyrene	12.63	202	92472	3.949	ng	98
81) Benzo(a)anthracene	13.82	228	46765	2.418	ng	98
83) Chrysene	13.85	228	42208	2.148	ng	98
88) Benzo(b)fluoranthene	14.84	252	48413m	2.083	ng	

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

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