

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062818\
 Data File : BF106950.D
 Acq On : 29 Jun 2018 4:29
 Operator : JU/SJ
 Sample : J3747-01
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RICH-REC-ROAD-STONE-4

Quant Time: Jun 29 07:03:19 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 25 18:50:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	62430	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	215074	20.00	ng	-0.01
38) Acenaphthene-d10	10.00	164	88558	20.00	ng	-0.01
63) Phenanthrene-d10	11.48	188	158127	20.00	ng	-0.02
75) Chrysene-d12	14.14	240	206609	20.00	ng	-0.03
86) Perylene-d12	15.67	264	170403	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	386407	104.81	ng	0.00
7) Phenol-d6	6.60	99	466626	101.35	ng	-0.02
23) Nitrobenzene-d5	7.52	82	286243	73.96	ng	-0.02
41) 2,4,6-Tribromophenol	10.79	330	109243	106.43	ng	-0.01
44) 2-Fluorobiphenyl	9.31	172	437711	83.38	ng	-0.02
78) Terphenyl-d14	13.07	244	531574	62.32	ng	-0.02
Target Compounds						
49) Dimethylphthalate	9.70	163	91337	13.752	ng	98
70) Phenanthrene	11.51	178	18510	2.427	ng	99
74) Fluoranthene	12.70	202	47632	5.666	ng	96
77) Pyrene	12.94	202	46119	3.385	ng	98
80) Benzo(a)anthracene	14.12	228	29362	2.332	ng	97
82) Chrysene	14.16	228	25965	2.241	ng	97
83) Bis(2-ethylhexyl)phthalate	14.09	149	22547	3.148	ng	# 97
87) Benzo(b)fluoranthene	15.21	252	39112	3.695	ng	# 92
88) Benzo(k)fluoranthene	15.21	252	39112	3.880	ng	93
89) Benzo(a)pyrene	15.60	252	19972	2.122	ng	99

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

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