

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061218\
 Data File : BF106340.D
 Acq On : 12 Jun 2018 13:21
 Operator : JU/SJ
 Sample : J3413-02
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 D-5

Quant Time: Jun 13 01:37:15 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 11 15:00:22 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	170407	20.00	ng	0.00
21) Naphthalene-d8	8.26	136	573653	20.00	ng	0.00
38) Acenaphthene-d10	10.02	164	254631	20.00	ng	0.01
63) Phenanthrene-d10	11.51	188	449597	20.00	ng	0.00
75) Chrysene-d12	14.15	240	498110	20.00	ng	0.00
86) Perylene-d12	15.70	264	517576	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	1320073	131.11	ng	0.02
7) Phenol-d6	6.60	99	1647722	129.53	ng	0.01
23) Nitrobenzene-d5	7.54	82	1052652	107.95	ng	0.00
41) 2,4,6-Tribromophenol	10.82	330	397292	162.16	ng	0.01
44) 2-Fluorobiphenyl	9.34	172	1408366	105.93	ng	0.00
78) Terphenyl-d14	13.10	244	1499907	74.79	ng	0.00
Target Compounds						
10) Phenol	6.61	94	34832	2.508	ng	97
37) 2-Methylnaphthalene	8.98	142	1735171	98.241	ng	# 94
49) Dimethylphthalate	9.73	163	196880	10.822	ng	# 91
57) Fluorene	10.58	166	273884	16.454	ng	# 83
70) Phenanthrene	11.54	178	736119	33.432	ng	96
77) Pyrene	12.95	202	155650	4.991	ng	# 96

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

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