

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121618\
 Data File : BF111526.D
 Acq On : 16 Dec 2018 23:42
 Operator : JU/SJ
 Sample : J6377-04
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P001-SD014-0612-01

Quant Time: Dec 18 12:11:39 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 14 13:26:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	122260	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	472118	20.00	ng	-0.01
39) Acenaphthene-d10	9.82	164	190846	20.00	ng	-0.01
64) Phenanthrene-d10	11.30	188	398435	20.00	ng	-0.01
76) Chrysene-d12	13.94	240	375964	20.00	ng	0.00
87) Perylene-d12	15.38	264	275078	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	1083685	147.50	ng	0.00
7) Phenol-d6	6.44	99	1111280	113.71	ng	0.00
23) Nitrobenzene-d5	7.35	82	741050	93.47	ng	0.00
42) 2,4,6-Tribromophenol	10.61	330	179703	105.12	ng	-0.01
45) 2-Fluorobiphenyl	9.15	172	1316891	106.58	ng	-0.01
79) Terphenyl-d14	12.89	244	1295639	80.92	ng	0.00
Target Compounds						
10) Phenol	6.45	94	71053	6.527	ng	89
13) 1,4-Dichlorobenzene	6.80	146	145825	14.867	ng	98
37) 2-Methylnaphthalene	8.78	142	31417	2.103	ng	97
50) Dimethylphthalate	9.54	163	70227	5.110	ng	# 98
71) Phenanthrene	11.33	178	46109	2.246	ng	94

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

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