

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122718\
 Data File : BF111767.D
 Acq On : 27 Dec 2018 17:25
 Operator : JU/SJ
 Sample : J6426-19 2X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P001-SS020-0002-01

Quant Time: Dec 28 06:56:17 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 19 17:39:39 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.91	152	135850	20.00	ng	0.01
21) Naphthalene-d8	8.19	136	460651	20.00	ng	0.01
39) Acenaphthene-d10	9.95	164	208363	20.00	ng	0.01
64) Phenanthrene-d10	11.43	188	377078	20.00	ng	0.02
76) Chrysene-d12	14.08	240	317602	20.00	ng	0.03
87) Perylene-d12	15.59	264	228678	20.00	ng	0.07

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.55	112	366107	45.94	ng	0.03
7) Phenol-d6	6.55	99	435281	42.91	ng	0.03
23) Nitrobenzene-d5	7.46	82	270522	34.18	ng	0.00
42) 2,4,6-Tribromophenol	10.74	330	113661	46.17	ng	0.02
45) 2-Fluorobiphenyl	9.26	172	497926	34.83	ng	0.00
79) Terphenyl-d14	13.02	244	526009	31.41	ng	0.02

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
9) Benzaldehyde	6.45	77	89451	12.823	ng	96
10) Phenol	6.56	94	69287	6.054	ng	98
17) 2-Methylphenol	7.17	107	16436	2.325	ng	94
20) 3+4-Methylphenols	7.32	107	96595	10.814	ng	# 63
27) 2,4-Dimethylphenol	7.85	122	95290	14.370	ng	88
31) Naphthalene	8.21	128	65457	2.957	ng	97
37) 2-Methylnaphthalene	8.90	142	296314	20.577	ng	99
38) 1-Methylnaphthalene	9.00	142	236492	17.100	ng	98
46) 1,1'-Biphenyl	9.36	154	54104	3.076	ng	95
50) Dimethylphthalate	9.65	163	65269	4.284	ng	# 99

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

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