

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122718\
 Data File : BF111768.D
 Acq On : 27 Dec 2018 17:51
 Operator : JU/SJ
 Sample : J6426-15 2X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P001-SS013C-1824-01

Quant Time: Dec 28 07:01:52 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	117833	20.00	ng	0.01
21) Naphthalene-d8	8.19	136	397545	20.00	ng	0.01
39) Acenaphthene-d10	9.95	164	178832	20.00	ng	0.02
64) Phenanthrene-d10	11.45	188	306253	20.00	ng	0.03
76) Chrysene-d12	14.11	240	199245	20.00	ng	0.06
87) Perylene-d12	15.62	264	168343	20.00	ng	0.11

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.55	112	109415	15.83	ng	0.04
7) Phenol-d6	6.55	99	181689	20.65	ng	0.03
23) Nitrobenzene-d5	7.46	82	143578	21.02	ng	0.00
42) 2,4,6-Tribromophenol	10.75	330	50490	23.89	ng	0.02
45) 2-Fluorobiphenyl	9.26	172	232039	18.91	ng	0.00
79) Terphenyl-d14	13.04	244	200814	19.11	ng	0.04

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
10) Phenol	6.57	94	313614	31.593	ng	97
17) 2-Methylphenol	7.17	107	20478	3.340	ng	# 82
20) 3+4-Methylphenols	7.32	107	102567	13.238	ng	# 64
30) 1,2,4-Trichlorobenzene	8.13	180	128484	18.731	ng	96
31) Naphthalene	8.22	128	173104	9.062	ng	97
37) 2-Methylnaphthalene	8.90	142	183094	14.733	ng	96
38) 1-Methylnaphthalene	9.00	142	104704	8.772	ng	# 98
40) 1,2,4,5-Tetrachlorobenzene	9.07	216	27332	4.651	ng	# 85
46) 1,1'-Biphenyl	9.37	154	38628	2.559	ng	97
50) Dimethylphthalate	9.66	163	37203	2.845	ng	# 60
71) Phenanthrene	11.47	178	88399	5.443	ng	# 85
78) Pvrene	12.90	202	95263	6.771	ng	# 83

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

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