

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF010719\
 Data File : BF111880.D
 Acq On : 7 Jan 2019 16:59
 Operator : JU/SJ
 Sample : K1053-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMP-1-DRUM

Quant Time: Jan 08 05:55:51 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF010719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 07 14:33:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	131040	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	550298	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	219457	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	383122	20.00	ng	0.00
76) Chrysene-d12	14.05	240	335175	20.00	ng	0.00
87) Perylene-d12	15.53	264	431216	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.54	112	403954	52.74	ng	0.02
7) Phenol-d6	6.55	99	293012	29.68	ng	0.01
23) Nitrobenzene-d5	7.46	82	727585	70.66	ng	0.00
42) 2,4,6-Tribromophenol	10.72	330	314300	110.11	ng	0.00
45) 2-Fluorobiphenyl	9.26	172	1308739	86.96	ng	0.00
79) Terphenyl-d14	13.00	244	1308344	74.10	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.69	88	20406	6.093	ng	# 54
9) Benzaldehyde	6.44	77	35406	4.177	ng	# 99
10) Phenol	6.56	94	1489231	138.794	ng	# 77
17) 2-Methylphenol	7.16	107	972403	142.255	ng	# 99
20) 3+4-Methylphenols	7.35	107	5777285	679.093	ng	# 68
27) 2,4-Dimethylphenol	7.86	122	4264662	575.122	ng	# 97
31) Naphthalene	8.20	128	1177466	45.074	ng	# 100
32) Benzoic acid	8.02	122	6309511	1311.192	ng	# 15
37) 2-Methylnaphthalene	8.89	142	155678	9.618	ng	# 94
38) 1-Methylnaphthalene	8.99	142	98670	6.355	ng	# 98
66) n-Nitrosodiphenylamine	10.58	169	66378	5.163	ng	# 96

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

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