

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060719\
 Data File : BF114907.D
 Acq On : 7 Jun 2019 23:12
 Operator : HP/JU
 Sample : K3214-03
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-4A

Quant Time: Jun 10 07:43:54 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF060419.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 04 16:57:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.66	152	335510	20.00	ng	0.00
21) Naphthalene-d8	7.95	136	880760	20.00	ng	0.00
39) Acenaphthene-d10	9.70	164	589233	20.00	ng	0.00
64) Phenanthrene-d10	11.16	188	1139836	20.00	ng	0.00
76) Chrysene-d12	13.78	240	681921	20.00	ng	-0.01
87) Perylene-d12	15.16	264	710941	20.00	ng	-0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.29	112	1539946m	80.20	ng	0.03
7) Phenol-d6	6.32	99	1192453	51.04	ng	0.02
23) Nitrobenzene-d5	7.23	82	1826767	127.68	ng	0.00
42) 2,4,6-Tribromophenol	10.48	330	815443	128.42	ng	0.00
45) 2-Fluorobiphenyl	9.03	172	3156693	100.36	ng	0.00
79) Terphenyl-d14	12.75	244	3227841	92.05	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.32	88	21717	2.369	ng	98
20) 3+4-Methylphenols	7.11	107	1232791	58.940	ng	# 30
31) Naphthalene	7.99	128	16882094	425.756	ng	# 93
37) 2-Methylnaphthalene	8.67	142	4850887	176.375	ng	# 94
38) 1-Methylnaphthalene	8.78	142	5821777	223.208	ng	# 99
46) 1,1'-Biphenyl	9.13	154	1194947	27.362	ng	96
52) Acenaphthene	9.73	154	2184275	61.963	ng	99
55) Dibenzofuran	9.90	168	2083593	42.506	ng	99
58) Fluorene	10.24	166	1464143	47.017	ng	100
71) Phenanthrene	11.19	178	1555915	28.594	ng	98
72) Anthracene	11.24	178	327048	5.711	ng	97
73) Carbazole	11.40	167	227598	4.636	ng	96
75) Fluoranthene	12.36	202	166498	2.940	ng	98

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

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