

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080918\
 Data File : BF108050.D
 Acq On : 9 Aug 2018 14:13
 Operator : JU/SJ
 Sample : J4262-07DL 2X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DUP-073118DL

Quant Time: Aug 09 17:57:20 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF080818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 08 12:24:24 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.01	152	330911	20.00	ng	0.00
21) Naphthalene-d8	8.30	136	1231758	20.00	ng	0.00
38) Acenaphthene-d10	10.07	164	593910	20.00	ng	0.00
63) Phenanthrene-d10	11.56	188	972392	20.00	ng	0.00
75) Chrysene-d12	14.21	240	780757	20.00	ng	0.00
86) Perylene-d12	15.78	264	655710	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.63	112	566322	30.98	ng	0.00
7) Phenol-d6	6.62	99	482952	19.88	ng	-0.01
23) Nitrobenzene-d5	7.57	82	1020924	46.25	ng	0.00
41) 2,4,6-Tribromophenol	10.85	330	346305	58.46	ng	0.00
44) 2-Fluorobiphenyl	9.37	172	1823556	49.29	ng	0.00
78) Terphenyl-d14	13.15	244	1522050	49.57	ng	0.00

Target Compounds				Qvalue		
17) 2-Methylphenol	7.24	107	43346	2.455	ng	95
20) 3+4-Methylphenols	7.40	107	97403	4.306	ng	88
27) 2,4-Dimethylphenol	7.93	122	74658	3.998	ng	96
31) Naphthalene	8.32	128	2950649	52.673	ng	# 92
37) 2-Methylnaphthalene	9.01	142	381772	9.633	ng	97
48) Acenaphthylene	9.92	152	191799	3.505	ng	98
49) Dimethylphthalate	9.77	163	94291	2.279	ng	# 98
51) Acenaphthene	10.10	154	609420	18.185	ng	99
54) Dibenzofuran	10.27	168	402737	8.295	ng	98
57) Fluorene	10.61	166	461339	12.003	ng	100
70) Phenanthrene	11.58	178	503515	10.978	ng	99
71) Anthracene	11.64	178	120600	2.566	ng	98
72) Carbazole	11.79	167	684919	16.399	ng	98
74) Fluoranthene	12.78	202	104217	2.082	ng	93
77) Pyrene	13.01	202	117211	2.371	ng	99

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

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