

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122118\
 Data File : BM018352.D
 Acq On : 21 Dec 2018 10:02
 Operator : JU/SJ
 Sample : J6347-05DL 5X
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Dec 21 12:40:32 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM121518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Dec 15 21:39:37 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.49	152	112853	20.00	ng	-0.02
21) Naphthalene-d8	10.25	136	466723	20.00	ng	-0.02
39) Acenaphthene-d10	14.15	164	283387	20.00	ng	-0.02
64) Phenanthrene-d10	16.91	188	656551	20.00	ng	-0.01
76) Chrysene-d12	21.14	240	663522	20.00	ng	0.00
87) Perylene-d12	23.30	264	553931	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.13	112	249484	36.93	ng	0.00
7) Phenol-d6	6.69	99	330230	35.17	ng	-0.01
23) Nitrobenzene-d5	8.65	82	239242	26.29	ng	-0.02
42) 2,4,6-Tribromophenol	15.66	330	138408	33.28	ng	0.00
45) 2-Fluorobiphenyl	12.75	172	433628	19.82	ng	-0.02
79) Terphenyl-d14	19.57	244	689258	23.49	ng	0.00

Target Compounds				Qvalue		
9) Benzaldehyde	6.63	77	26320	4.599	ng	# 1
10) Phenol	6.72	94	37249	3.746	ng	# 68
13) 1,4-Dichlorobenzene	7.52	146	156157	17.253	ng	98
14) 1,2-Dichlorobenzene	7.83	146	51085	5.855	ng	98
18) Hexachloroethane	8.57	117	32544	9.810	ng	# 1
30) 1,2,4-Trichlorobenzene	10.12	180	151880	18.739	ng	96
31) Naphthalene	10.30	128	670927	29.399	ng	99
35) Caprolactam	11.26	113	16554	6.097	ng	# 7
37) 2-Methylnaphthalene	11.93	142	466703	28.998	ng	96
38) 1-Methylnaphthalene	12.15	142	250848	15.973	ng	# 99
46) 1,1'-Biphenyl	12.96	154	160249	6.286	ng	98
48) 2-Nitroaniline	13.19	65	19847	3.434	ng	# 12
69) Atrazine	16.37	200	14923	2.036	ng	# 51
71) Phenanthrene	16.96	178	235119	6.593	ng	98

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

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