

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122118\
 Data File : BM018365.D
 Acq On : 21 Dec 2018 20:03
 Operator : JU/SJ
 Sample : J6389-06
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P001-SS013-0612-01

Quant Time: Dec 22 05:08:08 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	7.48	152	101048	20.00	ng	-0.02
21) Naphthalene-d8	10.25	136	362682	20.00	ng	-0.02
39) Acenaphthene-d10	14.15	164	184924	20.00	ng	-0.02
64) Phenanthrene-d10	16.91	188	374217	20.00	ng	-0.01
76) Chrysene-d12	21.14	240	438758	20.00	ng	0.00
87) Perylene-d12	23.31	264	471177	20.00	ng	0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.13	112	693701	114.68	ng	0.00
7) Phenol-d6	6.70	99	887757	105.59	ng	0.00
23) Nitrobenzene-d5	8.64	82	597423	84.49	ng	-0.02
42) 2,4,6-Tribromophenol	15.66	330	275076	101.35	ng	-0.01
45) 2-Fluorobiphenyl	12.75	172	1077420	75.46	ng	-0.02
79) Terphenyl-d14	19.57	244	1335716	68.85	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
9) Benzaldehyde	6.65	77	33504	6.538	ng	# 79
10) Phenol	6.73	94	101709	11.423	ng	92
20) 3+4-Methylphenols	8.29	107	48792	5.857	ng	94
27) 2,4-Dimethylphenol	9.47	122	20993	4.204	ng	95
49) Acenaphthylene	13.86	152	38185	2.069	ng	# 94
50) Dimethylphthalate	13.62	163	99748	6.488	ng	97
71) Phenanthrene	16.95	178	109217	5.373	ng	99
75) Fluoranthene	18.99	202	67965	2.876	ng	92
78) Pyrene	19.36	202	122405	4.682	ng	97
83) Chrysene	21.18	228	56348	2.286	ng	# 80
90) Benzo(a)pyrene	23.22	252	55815	2.217	ng	# 67
92) Benzo(a,h,i)perylene	26.11	276	45420	2.061	ng	# 92

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

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