

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122118\
 Data File : BM018357.D
 Acq On : 21 Dec 2018 13:07
 Operator : JU/SJ
 Sample : J6389-13
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P001-SS015-0006-01

Quant Time: Dec 22 04:08:42 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	119183	20.00	ng	-0.02
21) Naphthalene-d8	10.25	136	428123	20.00	ng	-0.02
39) Acenaphthene-d10	14.14	164	207293	20.00	ng	-0.02
64) Phenanthrene-d10	16.91	188	374716	20.00	ng	-0.01
76) Chrysene-d12	21.14	240	441116	20.00	ng	0.00
87) Perylene-d12	23.31	264	455087	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.13	112	838628	117.54	ng	0.00
7) Phenol-d6	6.70	99	1094240	110.34	ng	0.00
23) Nitrobenzene-d5	8.65	82	730693	87.54	ng	-0.02
42) 2,4,6-Tribromophenol	15.66	330	276136	90.76	ng	-0.01
45) 2-Fluorobiphenyl	12.75	172	1232959	77.04	ng	-0.02
79) Terphenyl-d14	19.57	244	1272977	65.27	ng	0.00
Target Compounds						
10) Phenol	6.73	94	95822	9.124	ng	91
40) 1,2,4,5-Tetrachlorobenzene	12.31	216	16787	2.355	ng	95
46) 1,1'-Biphenyl	12.97	154	58195	3.121	ng	92
50) Dimethylphthalate	13.62	163	117533	6.820	ng	96
68) Hexachlorobenzene	16.22	284	40218	8.089	ng	95

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

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