

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM121718\
 Data File : BM018241.D
 Acq On : 17 Dec 2018 10:19
 Operator : JU/SJ
 Sample : J6347-20
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P001-SD001-0006-01

Quant Time: Dec 18 00:48:43 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	128611	20.00	ng	-0.01
21) Naphthalene-d8	10.26	136	516026	20.00	ng	0.00
39) Acenaphthene-d10	14.15	164	299857	20.00	ng	-0.01
64) Phenanthrene-d10	16.92	188	675739	20.00	ng	0.00
76) Chrysene-d12	21.13	240	700534	20.00	ng	-0.01
87) Perylene-d12	23.30	264	620061	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.13	112	1350310	175.39	ng	0.00
7) Phenol-d6	6.70	99	1839413	171.89	ng	0.00
23) Nitrobenzene-d5	8.66	82	1065138	105.88	ng	0.00
42) 2,4,6-Tribromophenol	15.66	330	672743	152.86	ng	0.00
45) 2-Fluorobiphenyl	12.76	172	2173578	93.89	ng	0.00
79) Terphenyl-d14	19.57	244	2940505	94.93	ng	0.00
Target Compounds						
10) Phenol	6.72	94	118129	10.424	ng	97
13) 1,4-Dichlorobenzene	7.53	146	41886	4.061	ng	98
50) Dimethylphthalate	13.63	163	174723	7.009	ng	99

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

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