

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122118\
 Data File : BM018376.D
 Acq On : 22 Dec 2018 02:43
 Operator : JU/SJ
 Sample : J6389-09
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P001-SS014-0006-01

Quant Time: Dec 22 05:59:34 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	115235	20.00	ng	-0.02
21) Naphthalene-d8	10.25	136	465914	20.00	ng	-0.02
39) Acenaphthene-d10	14.14	164	249158	20.00	ng	-0.02
64) Phenanthrene-d10	16.91	188	465726	20.00	ng	-0.01
76) Chrysene-d12	21.13	240	419579	20.00	ng	-0.01
87) Perylene-d12	23.30	264	453572	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.12	112	719656	104.32	ng	0.00
7) Phenol-d6	6.70	99	984435	102.67	ng	0.00
23) Nitrobenzene-d5	8.64	82	669679	73.73	ng	-0.02
42) 2,4,6-Tribromophenol	15.66	330	317000	86.68	ng	-0.01
45) 2-Fluorobiphenyl	12.75	172	1248925	64.92	ng	-0.02
79) Terphenyl-d14	19.57	244	1130862	60.96	ng	0.00
Target Compounds						
6) Aniline	6.83	93	38619	3.211	ng	Qvalue # 86
10) Phenol	6.72	94	132398	13.039	ng	91
20) 3+4-Methylphenols	8.28	107	24177	2.545	ng	89
50) Dimethylphthalate	13.62	163	76978	3.716	ng	# 1

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

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