

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121818\
 Data File : BM018282.D
 Acq On : 18 Dec 2018 21:48
 Operator : JU/SJ
 Sample : J6340-11
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SB-10A

Quant Time: Dec 19 03:49:17 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	147093	20.00	ng	-0.02
21) Naphthalene-d8	10.25	136	546827	20.00	ng	-0.02
39) Acenaphthene-d10	14.14	164	266338	20.00	ng	-0.02
64) Phenanthrene-d10	16.91	188	461432	20.00	ng	-0.01
76) Chrysene-d12	21.13	240	456051	20.00	ng	-0.01
87) Perylene-d12	23.29	264	525926	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.12	112	1303329	148.01	ng	0.00
7) Phenol-d6	6.69	99	1757580	143.61	ng	-0.01
23) Nitrobenzene-d5	8.65	82	1117385	104.81	ng	-0.02
42) 2,4,6-Tribromophenol	15.66	330	429355	109.83	ng	-0.01
45) 2-Fluorobiphenyl	12.75	172	2049170	99.65	ng	-0.02
79) Terphenyl-d14	19.56	244	1824991	90.50	ng	-0.01
Target Compounds						
10) Phenol	6.72	94	82330	6.352	ng	93
50) Dimethylphthalate	13.62	163	206705	9.335	ng	97
71) Phenanthrene	16.95	178	117083	4.671	ng	99
75) Fluoranthene	18.98	202	196355	6.738	ng	99
78) Pyrene	19.35	202	179159	6.593	ng	99
81) Benzo(a)anthracene	21.12	228	103376	3.880	ng	99
83) Chrysene	21.17	228	98315	3.837	ng	97
88) Benzo(b)fluoranthene	22.66	252	103175	3.492	ng	93
90) Benzo(a)pyrene	23.20	252	63432	2.258	ng	94

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

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