

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM122816\
 Data File : BM008702.D
 Acq On : 28 Dec 2016 20:20
 Operator : UM/SJ
 Sample : H6067-20
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DA7H6

Manual Integrations
 APPROVED

mohammad
 12/30/2016 8:53:47 AM

Quant Time: Dec 29 04:30:51 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM121916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 29 03:29:39 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	418m	0.40	ng/ul	-0.17
2) Naphthalene-d8	10.43	136	1409	0.40	ng/ul	-0.13
6) Acenaphthene-d10	14.28	164	983m	0.40	ng/ul	-0.07
10) Phenanthrene-d10	17.04	188	1827m	0.40	ng/ul	-0.07
16) Chrysene-d12	21.28	240	1730	0.40	ng/ul	0.00
20) Perylene-d12	23.51	264	1456m	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.06	152	439	0.20	ng/ul	-0.10
14) Fluoranthene-d10	19.09	212	1162	0.23	ng/ul	-0.04
Target Compounds						
3) Naphthalene	10.48	128	431	0.11	ng/ul#	65
5) 2-Methylnaphthalene	12.13	142	515	0.19	ng/ul	100
12) Phenanthrene	17.07	178	1303m	0.23	ng/ul	
15) Fluoranthene	19.11	202	813	0.12	ng/ul	77
17) Pyrene	19.48	202	837	0.13	ng/ul#	65
18) Benzo(a)anthracene	21.26	228	463	0.08	ng/ul#	57
19) Chrysene	21.31	228	450	0.07	ng/ul#	61
23) Benzo(a)pyrene	23.38	252	307	0.06	ng/ul#	10
24) Indeno(1,2,3-cd)pyrene	25.79	276	206	0.03	ng/ul#	92

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

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