

Data Path : Z:\HPCHEM1\BNA_M\Data\BM120216\
 Data File : BM008158.D
 Acq On : 02 Dec 2016 16:15
 Operator : UM/SJ
 Sample : H5730-11MSD
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4WH6MSD

Quant Time: Dec 02 17:58:25 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM112116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 02 15:28:31 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	598	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.48	136	2109	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.33	164	1314	0.40	ng/ul	-0.02
10) Phenanthrene-d10	17.18	188	136356	0.40	ng/ul	0.07
16) Chrysene-d12	21.28	240	1955	0.40	ng/ul	-0.01
20) Perylene-d12	23.37	264	99117	0.40	ng/ul	-0.16
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.08	152	831	0.28	ng/ul	-0.01
14) Fluoranthene-d10	19.12	212	1647	0.00	ng/ul	-0.02
Target Compounds						
3) Naphthalene	10.52	128	208	0.04	ng/ul#	43
5) 2-Methylnaphthalene	12.15	142	85	0.02	ng/ul	100
7) Acenaphthylene	14.04	152	627	0.11	ng/ul#	81
8) Acenaphthene	14.38	153	1602	0.35	ng/ul	91
9) Fluorene	15.39	166	522	0.10	ng/ul#	96
15) Fluoranthene	19.14	202	15506	0.03	ng/ul	94
17) Pyrene	19.51	202	15904	2.34	ng/ul	98
18) Benzo(a)anthracene	21.26	228	5386	0.90	ng/ul	98
19) Chrysene	21.31	228	5487	0.74	ng/ul	96
22) Benzo(k)fluoranthene	22.85	252	9210	0.02	ng/ul#	94

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

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