

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM112116\
 Data File : BM007983.D
 Acq On : 22 Nov 2016 05:59
 Operator : UM/SJ
 Sample : H5694-14
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JHSS1

Manual Integrations
 APPROVED

sohil
 11/22/2016 4:00:28 PM

Quant Time: Nov 22 10:40:26 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM112116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 22 05:49:20 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	565	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.50	136	1872	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.35	164	1203	0.40	ng/ul	-0.02
10) Phenanthrene-d10	17.11	188	2656m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.30	240	2760	0.40	ng/ul	-0.01
20) Perylene-d12	23.54	264	2371m	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	601	0.23	ng/ul	-0.01
14) Fluoranthene-d10	19.14	212	1489	0.20	ng/ul	-0.01
Target Compounds						
3) Naphthalene	10.55	128	601	0.11	ng/ul#	78
5) 2-Methylnaphthalene	12.18	142	387	0.11	ng/ul	97
15) Fluoranthene	19.16	202	1896	0.18	ng/ul	95
17) Pyrene	19.52	202	2471	0.26	ng/ul	95
18) Benzo(a)anthracene	21.29	228	1196	0.14	ng/ul#	87
19) Chrysene	21.34	228	1569	0.15	ng/ul#	92
21) Benzo(b)fluoranthene	22.87	252	2135m	0.23	ng/ul	
22) Benzo(k)fluoranthene	22.90	252	766m	0.08	ng/ul	
23) Benzo(a)pyrene	23.43	252	1467m	0.17	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.80	276	1178m	0.11	ng/ul	
26) Benzo(g,h,i)perylene	26.50	276	1121m	0.12	ng/ul	

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

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