

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM120216\
 Data File : BM008177.D
 Acq On : 03 Dec 2016 03:35
 Operator : UM/SJ
 Sample : H5730-08DL 2X
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4WH4DL

Manual Integrations
 APPROVED

sohil
 12/5/2016 11:18:19 AM

Quant Time: Dec 03 05:05:42 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM112116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Dec 03 04:54:05 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.71	152	561	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.48	136	1851	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.33	164	1062	0.40	ng/ul	-0.02
10) Phenanthrene-d10	17.07	188	1859	0.40	ng/ul	-0.03
16) Chrysene-d12	21.29	240	1628	0.40	ng/ul	0.00
20) Perylene-d12	23.52	264	1674	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.09	152	312	0.12	ng/ul	-0.01
14) Fluoranthene-d10	19.13	212	691	0.13	ng/ul	0.00
Target Compounds						
					Qvalue	
9) Fluorene	15.39	166	111	0.03	ng/ul#	81
12) Phenanthrene	17.12	178	2285	0.39	ng/ul	93
13) Anthracene	17.21	178	459m	0.08	ng/ul	
15) Fluoranthene	19.15	202	4243	0.57	ng/ul	96
17) Pyrene	19.51	202	5493	0.97	ng/ul	97
18) Benzo(a)anthracene	21.27	228	2104	0.42	ng/ul	95
19) Chrysene	21.32	228	2587	0.42	ng/ul	95
21) Benzo(b)fluoranthene	22.85	252	2198m	0.33	ng/ul	
22) Benzo(k)fluoranthene	22.87	252	871m	0.13	ng/ul	
23) Benzo(a)pyrene	23.41	252	1914	0.31	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	25.78	276	1057	0.14	ng/ul#	90
26) Benzo(g,h,i)perylene	26.47	276	1031	0.15	ng/ul#	86

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

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