

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM120216\
 Data File : BM008154.D
 Acq On : 02 Dec 2016 13:52
 Operator : UM/SJ
 Sample : H5730-07
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4WH2

Manual Integrations
 APPROVED

sohil
 12/5/2016 11:17:25 AM

Quant Time: Dec 02 16:59:07 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	631	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.48	136	2094	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.33	164	1179	0.40	ng/ul	-0.02
10) Phenanthrene-d10	17.09	188	2086m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.28	240	1987	0.40	ng/ul	-0.01
20) Perylene-d12	23.52	264	1684m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.09	152	726	0.24	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	1366	0.23	ng/ul	-0.02
Target Compounds				Qvalue		
7) Acenaphthylene	14.04	152	846	0.16	ng/ul#	85
8) Acenaphthene	14.38	153	145	0.04	ng/ul#	85
9) Fluorene	15.39	166	290	0.06	ng/ul#	87
12) Phenanthrene	17.12	178	4311	0.66	ng/ul	94
13) Anthracene	17.21	178	1121	0.18	ng/ul#	86
15) Fluoranthene	19.14	202	8338m	0.99	ng/ul	
17) Pyrene	19.51	202	7964	1.15	ng/ul	98
18) Benzo(a)anthracene	21.26	228	3219	0.53	ng/ul	93
19) Chrysene	21.31	228	3884	0.51	ng/ul	95
21) Benzo(b)fluoranthene	22.85	252	5202m	0.77	ng/ul	
22) Benzo(k)fluoranthene	22.89	252	1179m	0.17	ng/ul	
23) Benzo(a)pyrene	23.42	252	3280m	0.52	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.77	276	2519m	0.32	ng/ul	
25) Dibenzo(a,h)anthracene	25.77	278	811m	0.13	ng/ul	
26) Benzo(g,h,i)perylene	26.45	276	2468m	0.36	ng/ul	

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

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