

Data Path : Z:\HPCHEM1\BNA M\DATA\BM112316\
 Data File : BM008054.D
 Acq On : 24 Nov 2016 03:12
 Operator : UM/SJ
 Sample : H5694-13
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JHSS0

Manual Integrations
 APPROVED

Umangi
 11/28/2016 5:28:27 PM

Quant Time: Nov 24 04:40:17 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM112116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 24 04:11:15 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	735	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.49	136	2469	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.35	164	1470	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.09	188	2977m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.29	240	2970	0.40	ng/ul	-0.01
20) Perylene-d12	23.53	264	2955	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.10	152	520	0.15	ng/ul	-0.02
14) Fluoranthene-d10	19.13	212	1290	0.15	ng/ul	-0.01
Target Compounds					Ovalue	
3) Naphthalene	10.54	128	1253	0.18	ng/ul#	92
5) 2-Methylnaphthalene	12.16	142	593	0.13	ng/ul	98
7) Acenaphthylene	14.06	152	814	0.12	ng/ul#	85
8) Acenaphthene	14.40	153	570	0.11	ng/ul	94
9) Fluorene	15.39	166	437	0.08	ng/ul#	71
12) Phenanthrene	17.12	178	10196	1.10	ng/ul	97
13) Anthracene	17.23	178	2850	0.33	ng/ul#	94
15) Fluoranthene	19.16	202	22369	1.87	ng/ul	94
17) Pyrene	19.52	202	25612	2.48	ng/ul	97
18) Benzo(a)anthracene	21.27	228	13081	1.43	ng/ul#	92
19) Chrysene	21.32	228	20288m	1.79	ng/ul	
21) Benzo(b)fluoranthene	22.86	252	26030m	2.21	ng/ul	
22) Benzo(k)fluoranthene	22.88	252	9956m	0.82	ng/ul	
23) Benzo(a)pyrene	23.43	252	18654	1.69	ng/ul	90
24) Indeno(1,2,3-cd)pyrene	25.78	276	15270	1.12	ng/ul#	93
25) Dibenzo(a,h)anthracene	25.79	278	4321	0.39	ng/ul#	56
26) Benzo(a,h,i)perylene	26.47	276	16231	1.34	ng/ul	97

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

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