

Data Path : Z:\HPCHEM1\BNA M\DATA\BM112316\
 Data File : BM008058.D
 Acq On : 24 Nov 2016 05:34
 Operator : UM/SJ
 Sample : H5694-21
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JHSS8

Manual Integrations
 APPROVED

Umangi
 11/28/2016 5:28:15 PM

Quant Time: Nov 24 06:26:01 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	674	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.49	136	2248	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.35	164	1334	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.09	188	2536	0.40	ng/ul	-0.02
16) Chrysene-d12	21.29	240	2577	0.40	ng/ul	-0.01
20) Perylene-d12	23.52	264	2594	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.10	152	412	0.13	ng/ul	-0.02
14) Fluoranthene-d10	19.13	212	1017	0.14	ng/ul	-0.01
Target Compounds					Ovalue	
3) Naphthalene	10.54	128	246	0.04	ng/ul#	45
5) 2-Methylnaphthalene	12.17	142	155	0.04	ng/ul	100
7) Acenaphthylene	14.06	152	298	0.05	ng/ul#	61
12) Phenanthrene	17.12	178	2227	0.28	ng/ul	96
13) Anthracene	17.23	178	488m	0.07	ng/ul	
15) Fluoranthene	19.16	202	5772	0.57	ng/ul	97
17) Pyrene	19.52	202	5805	0.65	ng/ul	99
18) Benzo(a)anthracene	21.27	228	3132	0.40	ng/ul#	91
19) Chrysene	21.32	228	3528	0.36	ng/ul	93
21) Benzo(b)fluoranthene	22.86	252	4691m	0.45	ng/ul	
22) Benzo(k)fluoranthene	22.89	252	1764m	0.17	ng/ul	
23) Benzo(a)pyrene	23.42	252	3340	0.34	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.78	276	2540	0.21	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.78	278	674	0.07	ng/ul#	21
26) Benzo(g,h,i)perylene	26.46	276	2668	0.25	ng/ul#	93

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

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