

Data Path : Z:\HPCHEM1\BNA M\DATA\BM112816\
 Data File : BM008092.D
 Acq On : 28 Nov 2016 18:32
 Operator : UM/SJ
 Sample : H5694-19DL 2X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JHSS6DL

Manual Integrations
 APPROVED

UMANGI
 11/29/2016 2:22:43 PM

Quant Time: Nov 29 09:48:03 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 29 00:11:14 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.71	152	661	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.49	136	2246	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.35	164	1322	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.09	188	2425	0.40	ng/ul	-0.02
16) Chrysene-d12	21.29	240	2499	0.40	ng/ul	0.00
20) Perylene-d12	23.53	264	2377	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.10	152	272	0.09	ng/ul	0.01
14) Fluoranthene-d10	19.13	212	640	0.09	ng/ul	-0.01
Target Compounds				Ovalue		
3) Naphthalene	10.54	128	420	0.07	ng/ul#	66
5) 2-Methylnaphthalene	12.16	142	223	0.06	ng/ul	98
7) Acenaphthylene	14.06	152	379	0.06	ng/ul#	69
8) Acenaphthene	14.40	153	149	0.03	ng/ul#	88
9) Fluorene	15.41	166	158	0.03	ng/ul#	82
12) Phenanthrene	17.14	178	3899	0.52	ng/ul	92
13) Anthracene	17.23	178	852m	0.12	ng/ul	
15) Fluoranthene	19.16	202	6967	0.71	ng/ul	98
17) Pyrene	19.52	202	6793	0.78	ng/ul	98
18) Benzo(a)anthracene	21.27	228	3147	0.41	ng/ul#	79
19) Chrysene	21.32	228	3108	0.33	ng/ul#	78
21) Benzo(b)fluoranthene	22.86	252	3990m	0.42	ng/ul	
22) Benzo(k)fluoranthene	22.90	252	1436m	0.15	ng/ul	
23) Benzo(a)pyrene	23.43	252	2995	0.34	ng/ul#	76
24) Indeno(1,2,3-cd)pyrene	25.79	276	2206	0.20	ng/ul#	90
26) Benzo(a,h,i)perylene	26.48	276	2194	0.23	ng/ul#	72

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

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