

Data Path : Z:\HPCHEM1\BNA_M\Data\BM112816\
 Data File : BM008090.D
 Acq On : 28 Nov 2016 17:22
 Operator : UM/SJ
 Sample : H5701-07DL 2X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4WD2DL

Quant Time: Nov 28 17:54:06 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 Nov 25 01:27:01 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.71	152	652	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.50	136	2277	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.35	164	1364	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.09	188	2540	0.40	ng/ul	-0.02
16) Chrysene-d12	21.30	240	2338	0.40	ng/ul	0.00
20) Perylene-d12	23.54	264	2358	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.10	152	257	0.08	ng/ul	0.00
14) Fluoranthene-d10	19.14	212	788	0.11	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
7) Acenaphthylene	14.06	152	841	0.14	ng/ul#	84
8) Acenaphthene	14.40	153	126	0.03	ng/ul#	85
9) Fluorene	15.41	166	199	0.04	ng/ul#	81
12) Phenanthrene	17.14	178	3136	0.40	ng/ul#	93
13) Anthracene	17.14	178	3134	0.42	ng/ul	95
15) Fluoranthene	19.16	202	9297	0.91	ng/ul	97
17) Pyrene	19.52	202	7956	0.98	ng/ul	99
18) Benzo(a)anthracene	21.28	228	5492	0.77	ng/ul#	80
19) Chrysene	21.33	228	5362	0.60	ng/ul#	78
21) Benzo(b)fluoranthene	22.86	252	10591	1.13	ng/ul#	75
22) Benzo(k)fluoranthene	22.86	252	10591	1.09	ng/ul#	73
23) Benzo(a)pyrene	23.43	252	5716	0.65	ng/ul#	75
24) Indeno(1,2,3-cd)pyrene	25.79	276	3876	0.36	ng/ul#	90
25) Dibenzo(a,h)anthracene	25.80	278	908	0.10	ng/ul#	1
26) Benzo(g,h,i)perylene	26.49	276	3253	0.34	ng/ul#	64

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

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