

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM032015\
 Data File : BM000740.D
 Acq On : 20 Mar 2015 14:11
 Operator : TP/IZ
 Sample : G1581-03
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 X1M13

Quant Time: Mar 23 11:10:57 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM031915.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 19 15:39:49 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	3198	0.40	ng/ul	0.00
2) Naphthalene-d8	10.41	136	11034	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.27	164	6032	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.02	188	10693	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	8354	0.40	ng/ul	0.01
20) Perylene-d12	23.42	264	7496	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.00	152	3904	0.31	ng/ul	0.00
14) Fluoranthene-d10	19.07	212	6545	0.31	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.46	128	10705	0.41	ng/ul	99
5) 2-Methylnaphthalene	12.08	142	14819	0.88	ng/ul	99
7) Acenaphthylene	14.00	152	11256	0.54	ng/ul	99
8) Acenaphthene	14.33	153	22523	1.33	ng/ul	99
12) Phenanthrene	17.05	178	16655	0.58	ng/ul	99
15) Fluoranthene	19.09	202	28902	0.91	ng/ul	99
17) Pyrene	19.45	202	23303	0.78	ng/ul	100
18) Benzo(a)anthracene	21.21	228	11794	0.51	ng/ul	94
21) Benzo(b)fluoranthene	22.76	252	18141	0.71	ng/ul	98
22) Benzo(k)fluoranthene	22.80	252	24417	0.93	ng/ul	97
23) Benzo(a)pyrene	23.31	252	11318	0.49	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.61	276	20227	0.75	ng/ul	98
25) Dibenzo(a,h)anthracene	25.62	278	20323	0.93	ng/ul	96
26) Benzo(g,h,i)perylene	26.27	276	22939	0.96	ng/ul	97

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

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