

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE072115\
 Data File : BE090179.D
 Acq On : 20 Jul 2015 17:33
 Operator : TP/IZ
 Sample : SSTD0.110
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Jul 21 05:11:52 2015

Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM02.2-EPA-SIM-BE072115.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Tue Jul 21 03:09:00 2015

Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	829	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	3999	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.21	164	2242	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	5906	0.40	ng/ul	0.00
16) Chrysene-d12	21.11	240	6227	0.40	ng/ul	0.00
20) Perylene-d12	23.43	264	5967	0.40	ng/ul	0.00

System Monitoring Compounds

4) 2-Methylnaphthalene-d10	11.96	152	629	0.10	ng/ul	0.00
14) Fluoranthene-d10	18.97	212	1740	0.10	ng/ul	0.00

Target Compounds

						Qvalue
3) Naphthalene	10.41	128	1065	0.10	ng/ul#	88
5) 2-Methylnaphthalene	12.04	142	725	0.10	ng/ul	95
7) Acenaphthylene	13.93	152	4764	0.10	ng/ul#	93
8) Acenaphthene	14.27	153	3458	0.10	ng/ul	95
9) Fluorene	15.27	166	1014	0.10	ng/ul#	97
12) Phenanthrene	16.99	178	7051	0.10	ng/ul	94
13) Anthracene	17.07	178	6545	0.09	ng/ul	94
15) Fluoranthene	19.00	202	7910	0.10	ng/ul	91
17) Pyrene	19.36	202	8370	0.10	ng/ul	95
18) Benzo(a)anthracene	21.10	228	1633	0.10	ng/ul	96
19) Chrysene	21.15	228	1987	0.09	ng/ul	95
21) Benzo(b)fluoranthene	22.72	252	1575	0.09	ng/ul	88
22) Benzo(k)fluoranthene	22.77	252	2053	0.08	ng/ul#	85
23) Benzo(a)pyrene	23.32	252	1602	0.09	ng/ul#	85
24) Indeno(1,2,3-cd)pyrene	25.82	276	7937	0.10	ng/ul	96
25) Dibenzo(a,h)anthracene	25.85	278	1458	0.10	ng/ul#	84
26) Benzo(g,h,i)perylene	26.56	276	6899	0.09	ng/ul#	89

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

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