

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM031915\
 Data File : BM000729.D
 Acq On : 19 Mar 2015 14:36
 Operator : TP/IZ
 Sample : SSTD1.699
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD1.699

Quant Time: Mar 19 15:18:36 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 14:57:23 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	2856	0.40	ng/ul	0.00
2) Naphthalene-d8	10.41	136	9724	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.27	164	5261	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.02	188	9442	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	8393	0.40	ng/ul	0.00
20) Perylene-d12	23.41	264	7326	0.40	ng/ul	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.00	152	18268	1.68	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	31686	1.80	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.46	128	38284	1.66	ng/ul	99
5) 2-Methylnaphthalene	12.08	142	24528	1.68	ng/ul	100
7) Acenaphthylene	14.00	152	31784	1.63	ng/ul	98
8) Acenaphthene	14.33	153	24868	1.55	ng/ul	100
9) Fluorene	15.34	166	27922	1.58	ng/ul	96
11) Pentachlorophenol	16.68	266	1703	2.00	ng/ul	98
12) Phenanthrene	17.05	178	42387	1.70	ng/ul	99
13) Anthracene	17.15	178	38503	1.79	ng/ul	98
15) Fluoranthene	19.09	202	47935	1.82	ng/ul	99
17) Pyrene	19.45	202	48469	1.56	ng/ul	99
18) Benzo(a)anthracene	21.20	228	39661	1.74	ng/ul	99
19) Chrysene	21.25	228	44641	1.62	ng/ul	99
21) Benzo(b)fluoranthene	22.75	252	41626	1.60	ng/ul	96
22) Benzo(k)fluoranthene	22.79	252	42108	1.61	ng/ul	95
23) Benzo(a)pyrene	23.31	252	38349	1.67	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	25.60	276	44978	1.70	ng/ul	100
25) Dibenzo(a,h)anthracene	25.61	278	36493	1.70	ng/ul	94
26) Benzo(g,h,i)perylene	26.26	276	39377	1.66	ng/ul	97

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

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