

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM031915\
 Data File : BM000727.D
 Acq On : 19 Mar 2015 13:25
 Operator : TP/IZ
 Sample : SSTD0.497
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.497

Quant Time: Mar 19 15:13:41 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	2782	0.40	ng/ul	0.00
2) Naphthalene-d8	10.41	136	9507	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.27	164	5126	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.02	188	9215	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	7808	0.40	ng/ul	0.00
20) Perylene-d12	23.42	264	7217	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.00	152	4973	0.47	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	8251	0.48	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.46	128	10345	0.46	ng/ul	100
5) 2-Methylnaphthalene	12.08	142	6543	0.46	ng/ul	100
7) Acenaphthylene	14.00	152	8108	0.43	ng/ul	100
8) Acenaphthene	14.33	153	6644	0.42	ng/ul	100
9) Fluorene	15.34	166	7294	0.42	ng/ul	100
11) Pentachlorophenol	16.68	266	435	0.52	ng/ul	100
12) Phenanthrene	17.05	178	11130	0.46	ng/ul	100
13) Anthracene	17.15	178	9785	0.47	ng/ul	100
15) Fluoranthene	19.09	202	12288	0.48	ng/ul	100
17) Pyrene	19.45	202	12621	0.44	ng/ul	100
18) Benzo(a)anthracene	21.20	228	9756	0.46	ng/ul	100
19) Chrysene	21.25	228	11776	0.46	ng/ul	100
21) Benzo(b)fluoranthene	22.75	252	11008	0.43	ng/ul	100
22) Benzo(k)fluoranthene	22.79	252	11957	0.46	ng/ul	100
23) Benzo(a)pyrene	23.31	252	10290	0.46	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.60	276	11976	0.46	ng/ul	100
25) Dibenzo(a,h)anthracene	25.62	278	9732	0.46	ng/ul	100
26) Benzo(g,h,i)perylene	26.26	276	10612	0.45	ng/ul	100

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

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