

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM082715\
 Data File : BM002555.D
 Acq On : 27 Aug 2015 15:01
 Operator : UM/IZ
 Sample : SSTD0.458
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.458

Quant Time: Aug 27 16:08:19 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM082715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 27 16:04:04 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.78	152	9168	0.40	ng/ul	0.00
2) Naphthalene-d8	11.64	136	37935	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.36	164	20076	0.40	ng/ul	0.00
10) Phenanthrene-d10	18.08	188	43739	0.40	ng/ul	0.00
16) Chrysene-d12	22.16	240	32661	0.40	ng/ul	0.00
20) Perylene-d12	24.78	264	24820	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	13.17	152	19388	0.33	ng/ul	0.00
14) Fluoranthene-d10	20.03	212	38778	0.33	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	11.69	128	36800	0.32	ng/ul	99
5) 2-Methylnaphthalene	13.25	142	25164	0.33	ng/ul	100
7) Acenaphthylene	15.09	152	39477	0.33	ng/ul	94
8) Acenaphthene	15.42	153	27800	0.33	ng/ul	85
9) Fluorene	16.39	166	32031	0.33	ng/ul	93
11) Pentachlorophenol	17.72	266	1364	0.12	ng/ul	96
12) Phenanthrene	18.11	178	50143	0.33	ng/ul	98
13) Anthracene	18.20	178	48364	0.33	ng/ul	98
15) Fluoranthene	20.07	202	52520	0.31	ng/ul	90
17) Pyrene	20.42	202	53325	0.36	ng/ul	97
18) Benzo(a)anthracene	22.14	228	39893	0.31	ng/ul	98
19) Chrysene	22.20	228	38289	0.30	ng/ul	93
21) Benzo(b)fluoranthene	23.96	252	34202	0.30	ng/ul	96
22) Benzo(k)fluoranthene	24.01	252	31049	0.33	ng/ul#	95
23) Benzo(a)pyrene	24.66	252	30478	0.31	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	27.54	276	35018	0.33	ng/ul	98
25) Dibenzo(a,h)anthracene	27.54	278	28561	0.34	ng/ul	99
26) Benzo(g,h,i)perylene	28.40	276	29707	0.33	ng/ul	98

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

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