

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM082715\
 Data File : BM002556.D
 Acq On : 27 Aug 2015 15:37
 Operator : UM/IZ
 Sample : SSTD0.859
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.859

Quant Time: Aug 27 16:30:35 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	9283	0.40	ng/ul	0.00
2) Naphthalene-d8	11.64	136	38396	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.36	164	20250	0.40	ng/ul	0.00
10) Phenanthrene-d10	18.08	188	43651	0.40	ng/ul	0.00
16) Chrysene-d12	22.16	240	30978	0.40	ng/ul	0.00
20) Perylene-d12	24.77	264	24195	0.40	ng/ul	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	13.17	152	42922	0.72	ng/ul	0.00
14) Fluoranthene-d10	20.03	212	81858	0.69	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	11.69	128	81245	0.70	ng/ul	98
5) 2-Methylnaphthalene	13.25	142	56127	0.73	ng/ul	100
7) Acenaphthylene	15.09	152	86865	0.73	ng/ul	94
8) Acenaphthene	15.42	153	61312	0.73	ng/ul	85
9) Fluorene	16.39	166	70887	0.73	ng/ul	94
11) Pentachlorophenol	17.72	266	3467	0.30	ng/ul	99
12) Phenanthrene	18.11	178	106300	0.70	ng/ul	98
13) Anthracene	18.20	178	104437	0.71	ng/ul	98
15) Fluoranthene	20.07	202	111693	0.66	ng/ul	89
17) Pyrene	20.42	202	111754	0.79	ng/ul	96
18) Benzo(a)anthracene	22.14	228	83343	0.69	ng/ul	96
19) Chrysene	22.19	228	78650	0.65	ng/ul	99
21) Benzo(b)fluoranthene	23.96	252	74916	0.67	ng/ul	98
22) Benzo(k)fluoranthene	24.01	252	63570	0.69	ng/ul	99
23) Benzo(a)pyrene	24.66	252	64835	0.68	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	27.53	276	74578	0.72	ng/ul	98
25) Dibenzo(a,h)anthracene	27.54	278	60974	0.74	ng/ul	97
26) Benzo(g,h,i)perylene	28.39	276	63672	0.73	ng/ul	98

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

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