

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM072816\
 Data File : BM006725.D
 Acq On : 28 Jul 2016 16:00
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.453

Quant Time: Jul 28 16:33:40 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM072816.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 28 15:44:03 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	183234	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.38	136	713646	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.26	164	402613	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.00	188	895031	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	790235	0.40	ng/ul	0.00
20) Perylene-d12	23.42	264	874446	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	11.99	152	6196	0.01	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	15785	0.01	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.43	128	251	0.00	ng/ul#	30
5) 2-Methylnaphthalene	12.06	142	81	0.00	ng/ul	100
7) Acenaphthylene	13.99	152	387	0.00	ng/ul#	43
8) Acenaphthene	14.31	153	117	0.00	ng/ul#	82
9) Fluorene	15.26	166	2063	0.00	ng/ul#	23
11) Pentachlorophenol	16.70	266	41	0.00	ng/ul#	81
12) Phenanthrene	17.04	178	1571	0.00	ng/ul#	76
13) Anthracene	17.11	178	750	0.00	ng/ul#	1
15) Fluoranthene	19.08	202	3004	0.00	ng/ul	88
17) Pyrene	19.44	202	3658	0.00	ng/ul#	60
18) Benzo(a)anthracene	21.21	228	5827	0.00	ng/ul#	81
19) Chrysene	21.21	228	5827	0.00	ng/ul#	78
21) Benzo(b)fluoranthene	22.76	252	2812	0.00	ng/ul#	47
22) Benzo(k)fluoranthene	22.76	252	2812	0.00	ng/ul#	49
23) Benzo(a)pyrene	23.27	252	4919	0.00	ng/ul#	47
24) Indeno(1,2,3-cd)pyrene	25.63	276	1315	0.00	ng/ul#	57
25) Dibenzo(a,h)anthracene	25.63	278	787	0.00	ng/ul#	5
26) Benzo(g,h,i)perylene	26.30	276	1009	0.00	ng/ul#	69

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

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