

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022820\
 Data File : BM025182.D
 Acq On : 29 Feb 2020 23:29
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 55 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.498

Quant Time: Mar 02 05:52:57 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM022420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 02 02:08:15 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.68	152	12153	0.40	ng/ul	0.00
2) Naphthalene-d8	10.45	136	50849	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	34793	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	83250	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	92341	0.40	ng/ul	0.00
20) Perylene-d12	23.43	264	93287	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.05	152	7557	0.08	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	18986	0.07	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.50	128	58941	0.381	ng/ul#	82
5) 2-Methylnaphthalene	12.12	142	43485	0.389	ng/ul	100
7) Acenaphthylene	14.03	152	72963	0.364	ng/ul#	93
8) Acenaphthene	14.38	153	50101	0.372	ng/ul	98
9) Fluorene	15.36	166	62135	0.367	ng/ul#	96
11) Pentachlorophenol	16.71	266	10940	0.430	ng/ul	99
12) Phenanthrene	17.09	178	104772	0.358	ng/ul	95
13) Anthracene	17.18	178	102749	0.362	ng/ul	95
15) Fluoranthene	19.11	202	130778	0.307	ng/ul	95
17) Pyrene	19.47	202	138808	0.334	ng/ul	95
18) Benzo(a)anthracene	21.22	228	140422	0.352	ng/ul	97
19) Chrysene	21.27	228	140593	0.354	ng/ul	97
21) Benzo(b)fluoranthene	22.78	252	143862	0.393	ng/ul	90
22) Benzo(k)fluoranthene	22.82	252	147718	0.408	ng/ul#	92
23) Benzo(a)pyrene	23.33	252	131592	0.394	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	25.61	276	168931	0.404	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.63	278	140761	0.400	ng/ul#	92
26) Benzo(g,h,i)perylene	26.28	276	140252	0.398	ng/ul#	90

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

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