

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022820\
 Data File : BM025168.D
 Acq On : 29 Feb 2020 15:06
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.497

Quant Time: Mar 02 01:18:57 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 01:15:39 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.68	152	2777	0.40	ng/ul	0.00
2) Naphthalene-d8	10.45	136	10818	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	6806	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	15682	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	16248	0.40	ng/ul	0.00
20) Perylene-d12	23.43	264	17875	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.05	152	7900	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	21380	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.50	128	13393	0.407	ng/ul#	83
5) 2-Methylnaphthalene	12.12	142	9287	0.391	ng/ul	100
7) Acenaphthylene	14.03	152	16473	0.420	ng/ul	95
8) Acenaphthene	14.37	153	10344	0.392	ng/ul	99
9) Fluorene	15.36	166	13240	0.400	ng/ul	95
11) Pentachlorophenol	16.71	266	2304	0.481	ng/ul	99
12) Phenanthrene	17.09	178	21537	0.390	ng/ul	97
13) Anthracene	17.18	178	21605	0.404	ng/ul	96
15) Fluoranthene	19.11	202	28216	0.352	ng/ul	98
17) Pyrene	19.47	202	30319	0.414	ng/ul	98
18) Benzo(a)anthracene	21.22	228	28612	0.408	ng/ul	99
19) Chrysene	21.27	228	27985	0.400	ng/ul	99
21) Benzo(b)fluoranthene	22.78	252	30032	0.429	ng/ul	95
22) Benzo(k)fluoranthene	22.82	252	29497	0.425	ng/ul	98
23) Benzo(a)pyrene	23.34	252	26887	0.420	ng/ul#	94
24) Indeno(1,2,3-cd)pyrene	25.62	276	34212	0.427	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.63	278	27801	0.413	ng/ul	98
26) Benzo(g,h,i)perylene	26.28	276	27885	0.413	ng/ul#	93

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

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