

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030320\
 Data File : BM025254.D
 Acq On : 04 Mar 2020 17:32
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.486

Quant Time: Mar 04 18:07:57 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM030320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 04 08:02:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.68	152	3755	0.40	ng/ul	0.00
2) Naphthalene-d8	10.45	136	15570	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	11005	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	26783	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	26304	0.40	ng/ul	0.00
20) Perylene-d12	23.43	264	23211	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	10358	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	30240	0.35	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.50	128	16573	0.354	ng/ul#	82
5) 2-Methylnaphthalene	12.11	142	12075	0.354	ng/ul	99
7) Acenaphthylene	14.02	152	18284	0.355	ng/ul	94
8) Acenaphthene	14.37	153	14205	0.355	ng/ul	99
9) Fluorene	15.36	166	17629	0.352	ng/ul	96
11) Pentachlorophenol	16.70	266	2581	0.368	ng/ul	98
12) Phenanthrene	17.09	178	30880	0.358	ng/ul	96
13) Anthracene	17.18	178	27383	0.357	ng/ul	95
15) Fluoranthene	19.11	202	37387	0.346	ng/ul	96
17) Pyrene	19.47	202	37398	0.375	ng/ul	96
18) Benzo(a)anthracene	21.22	228	35745	0.370	ng/ul	97
19) Chrysene	21.27	228	38303	0.361	ng/ul	97
21) Benzo(b)fluoranthene	22.78	252	34022	0.366	ng/ul	93
22) Benzo(k)fluoranthene	22.82	252	36320	0.373	ng/ul#	96
23) Benzo(a)pyrene	23.33	252	28668	0.368	ng/ul#	93
24) Indeno(1,2,3-cd)pyrene	25.61	276	32295	0.332	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.63	278	26803	0.333	ng/ul#	94
26) Benzo(g,h,i)perylene	26.27	276	25886	0.314	ng/ul#	91

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

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