

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030320\
 Data File : BM025234.D
 Acq On : 04 Mar 2020 04:28
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.484

Quant Time: Mar 04 04:59:49 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 00:52:51 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.68	152	3955	0.40	ng/ul	0.00
2) Naphthalene-d8	10.44	136	16545	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	11869	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	28972	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	28686	0.40	ng/ul	0.00
20) Perylene-d12	23.42	264	26289	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	11181	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	33018	0.35	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.49	128	17978	0.361	ng/ul#	83
5) 2-Methylnaphthalene	12.11	142	13063	0.360	ng/ul	98
7) Acenaphthylene	14.02	152	19441	0.350	ng/ul	94
8) Acenaphthene	14.37	153	15372	0.357	ng/ul	99
9) Fluorene	15.36	166	19006	0.352	ng/ul	96
11) Pentachlorophenol	16.70	266	2801	0.369	ng/ul	99
12) Phenanthrene	17.09	178	33336	0.357	ng/ul	95
13) Anthracene	17.18	178	29797	0.359	ng/ul	94
15) Fluoranthene	19.11	202	41228	0.353	ng/ul	95
17) Pyrene	19.47	202	40722	0.374	ng/ul	95
18) Benzo(a)anthracene	21.21	228	39272	0.372	ng/ul	98
19) Chrysene	21.27	228	42230	0.364	ng/ul	97
21) Benzo(b)fluoranthene	22.77	252	38572	0.367	ng/ul	92
22) Benzo(k)fluoranthene	22.81	252	40752	0.369	ng/ul#	96
23) Benzo(a)pyrene	23.32	252	32368	0.367	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.60	276	37482	0.341	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.62	278	31232	0.342	ng/ul#	94
26) Benzo(g,h,i)perylene	26.26	276	30151	0.323	ng/ul#	90

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

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