

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040521\
 Data File : BM029350.D
 Acq On : 05 Apr 2021 20:48
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.483

Quant Time: Apr 06 04:06:54 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM040521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 05 15:51:17 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	1978	0.40	ng/ul	0.00
2) Naphthalene-d8	10.48	136	6446	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.33	164	3183	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.07	188	6321	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	6413	0.40	ng/ul	0.00
20) Perylene-d12	23.46	264	6884	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.07	152	3553	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.10	212	6353	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.53	128	7341	0.421	ng/ul	99
5) 2-Methylnaphthalene	12.15	142	4600	0.407	ng/ul	100
7) Acenaphthylene	14.05	152	5345	0.396	ng/ul	100
8) Acenaphthene	14.39	153	4707	0.412	ng/ul	97
9) Fluorene	15.38	166	5101	0.403	ng/ul	99
11) Pentachlorophenol	16.73	266	513	0.358	ng/ul	97
12) Phenanthrene	17.11	178	7711	0.400	ng/ul	98
13) Anthracene	17.20	178	6603	0.390	ng/ul	99
15) Fluoranthene	19.13	202	7763	0.363	ng/ul	99
17) Pyrene	19.49	202	8030	0.388	ng/ul	99
18) Benzo(a)anthracene	21.23	228	6737	0.369	ng/ul	99
19) Chrysene	21.28	228	7770	0.386	ng/ul	99
21) Benzo(b)fluoranthene	22.80	252	8502	0.407	ng/ul	94
22) Benzo(k)fluoranthene	22.84	252	8766	0.410	ng/ul	98
23) Benzo(a)pyrene	23.36	252	7373	0.408	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	25.68	276	10278	0.410	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.70	278	8638	0.411	ng/ul	99
26) Benzo(g,h,i)perylene	26.36	276	8764	0.413	ng/ul	98

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

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