

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030620\
 Data File : BM025357.D
 Acq On : 07 Mar 2020 19:26
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.416

Quant Time: Mar 07 21:00:49 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 : Sat Mar 07 19:35:17 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	4249	0.40	ng/ul	0.00
2) Naphthalene-d8	10.44	136	18931	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	13781	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	32315	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	25105	0.40	ng/ul	0.00
20) Perylene-d12	23.41	264	21053	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	13132	0.37	ng/ul	-0.01
14) Fluoranthene-d10	19.07	212	34459	0.33	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.48	128	20711	0.364	ng/ul#	83
5) 2-Methylnaphthalene	12.10	142	15364	0.370	ng/ul	99
7) Acenaphthylene	14.01	152	22413	0.347	ng/ul	93
8) Acenaphthene	14.36	153	18208	0.364	ng/ul	99
9) Fluorene	15.35	166	22361	0.356	ng/ul	95
11) Pentachlorophenol	16.69	266	2579	0.305	ng/ul	98
12) Phenanthrene	17.08	178	37565	0.360	ng/ul	95
13) Anthracene	17.17	178	33904	0.367	ng/ul	95
15) Fluoranthene	19.10	202	43524	0.334	ng/ul	96
17) Pyrene	19.46	202	42456	0.446	ng/ul	97
18) Benzo(a)anthracene	21.21	228	35181	0.381	ng/ul	98
19) Chrysene	21.26	228	36644	0.361	ng/ul	97
21) Benzo(b)fluoranthene	22.76	252	30937	0.367	ng/ul	94
22) Benzo(k)fluoranthene	22.81	252	32538	0.368	ng/ul#	98
23) Benzo(a)pyrene	23.32	252	26250	0.371	ng/ul#	94
24) Indeno(1,2,3-cd)pyrene	25.59	276	32931	0.374	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.60	278	27242	0.373	ng/ul#	95
26) Benzo(g,h,i)perylene	26.25	276	26813	0.358	ng/ul#	91

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

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