

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080219\
 Data File : BM021777.D
 Acq On : 03 Aug 2019 06:35
 Operator : HP/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.421

Quant Time: Aug 03 07:31:54 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM080219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Aug 03 05:33:59 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	697	0.40	ng/ul	0.00
2) Naphthalene-d8	10.46	136	3400	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	1529	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.07	188	2642	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	1630	0.40	ng/ul	0.00
20) Perylene-d12	23.49	264	1678	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.07	152	1955	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.10	212	2368	0.39	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.50	128	3813	0.400	ng/ul	98
5) 2-Methylnaphthalene	12.14	142	2346	0.429	ng/ul	98
7) Acenaphthylene	14.03	152	2966	0.386	ng/ul	99
8) Acenaphthene	14.38	153	2564	0.394	ng/ul	99
9) Fluorene	15.38	166	2496	0.367	ng/ul	98
11) Pentachlorophenol	16.77	266	116	0.313	ng/ul#	87
12) Phenanthrene	17.12	178	3455	0.391	ng/ul	98
13) Anthracene	17.22	178	2741	0.407	ng/ul	99
15) Fluoranthene	19.13	202	3179	0.361	ng/ul	95
17) Pyrene	19.50	202	3388	0.470	ng/ul	96
18) Benzo(a)anthracene	21.25	228	2061	0.422	ng/ul	96
19) Chrysene	21.30	228	3106	0.379	ng/ul	99
21) Benzo(b)fluoranthene	22.82	252	2567	0.343	ng/ul	96
22) Benzo(k)fluoranthene	22.86	252	2788	0.354	ng/ul	97
23) Benzo(a)pyrene	23.39	252	2230	0.379	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.72	276	2645	0.389	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.75	278	2061	0.389	ng/ul	92
26) Benzo(g,h,i)perylene	26.41	276	2493	0.382	ng/ul	96

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

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