

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM030920\
 Data File : BM025414.D
 Acq On : 09 Mar 2020 12:58
 Operator : CG/JU
 Sample : SSTD3.278
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD3.278

Quant Time: Mar 09 13:29:32 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM030920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 09 12:03:47 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	2179	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	8675	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	6337	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	14562	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	12628	0.40	ng/ul	0.00
20) Perylene-d12	23.41	264	10170	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	52646	3.24	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	146950	3.10	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.48	128	82110	3.145	ng/ul	98
5) 2-Methylnaphthalene	12.10	142	62092	3.265	ng/ul	100
7) Acenaphthylene	14.01	152	86915	2.929	ng/ul	98
8) Acenaphthene	14.35	153	74188	3.223	ng/ul	99
9) Fluorene	15.34	166	94820	3.287	ng/ul	100
11) Pentachlorophenol	16.69	266	13621	3.574	ng/ul	99
12) Phenanthrene	17.07	178	154434	3.289	ng/ul	100
13) Anthracene	17.16	178	146967	3.527	ng/ul	99
15) Fluoranthene	19.09	202	186212	3.168	ng/ul	100
17) Pyrene	19.45	202	177865	3.714	ng/ul	100
18) Benzo(a)anthracene	21.20	228	153426	3.305	ng/ul	99
19) Chrysene	21.26	228	156602	3.070	ng/ul	98
21) Benzo(b)fluoranthene	22.75	252	138425	3.402	ng/ul	94
22) Benzo(k)fluoranthene	22.80	252	144165	3.378	ng/ul#	94
23) Benzo(a)pyrene	23.31	252	117865	3.450	ng/ul#	91
24) Indeno(1,2,3-cd)pyrene	25.58	276	153397	3.602	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.59	278	129946	3.683	ng/ul	95
26) Benzo(g,h,i)perylene	26.24	276	130060	3.597	ng/ul	98

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

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