

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM040521\
 Data File : BM029343.D
 Acq On : 05 Apr 2021 15:12
 Operator : CG/JU
 Sample : SSTD3.281
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD3.281

Quant Time: Apr 05 15:48:24 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 14:08:43 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	2105	0.40	ng/ul	0.00
2) Naphthalene-d8	10.48	136	7048	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.33	164	3680	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.07	188	7428	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	8393	0.40	ng/ul	0.00
20) Perylene-d12	23.45	264	8495	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.07	152	29732	2.97	ng/ul	0.00
14) Fluoranthene-d10	19.10	212	59940	3.13	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.53	128	58180	3.316	ng/ul	97
5) 2-Methylnaphthalene	12.15	142	38448	3.272	ng/ul	100
7) Acenaphthylene	14.05	152	49932	3.325	ng/ul	98
8) Acenaphthene	14.39	153	40651	3.414	ng/ul	100
9) Fluorene	15.38	166	45339	3.431	ng/ul	98
11) Pentachlorophenol	16.72	266	5679	2.926	ng/ul	95
12) Phenanthrene	17.11	178	69009	3.324	ng/ul	99
13) Anthracene	17.20	178	65012	3.284	ng/ul	98
15) Fluoranthene	19.13	202	75212	3.182	ng/ul	98
17) Pyrene	19.49	202	75768	2.485	ng/ul	98
18) Benzo(a)anthracene	21.23	228	68652	2.396	ng/ul	99
19) Chrysene	21.28	228	70926	2.490	ng/ul	99
21) Benzo(b)fluoranthene	22.79	252	77808	2.676	ng/ul	89
22) Benzo(k)fluoranthene	22.84	252	77761	2.680	ng/ul#	90
23) Benzo(a)pyrene	23.36	252	66987	2.587	ng/ul#	84
24) Indeno(1,2,3-cd)pyrene	25.68	276	93658	2.777	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.69	278	79997	2.827	ng/ul	94
26) Benzo(g,h,i)perylene	26.36	276	78772	2.806	ng/ul	96

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

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