

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM033020\
 Data File : BM025859.D
 Acq On : 30 Mar 2020 14:30
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SICV82

Quant Time: Mar 30 15:02:07 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM033020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 30 14:21:01 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	1957	0.40	ng/ul	0.00
2) Naphthalene-d8	10.33	136	7555	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	4390	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	10640	0.40	ng/ul	0.00
16) Chrysene-d12	21.15	240	9618	0.40	ng/ul	0.00
20) Perylene-d12	23.29	264	9136	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.93	152	5098	0.39	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	12490	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.38	128	8179	0.364	ng/ul	99
5) 2-Methylnaphthalene	12.00	142	5832	0.386	ng/ul	100
7) Acenaphthylene	13.92	152	7799	0.419	ng/ul	99
8) Acenaphthene	14.27	153	6112	0.393	ng/ul	99
9) Fluorene	15.26	166	7329	0.398	ng/ul	98
11) Pentachlorophenol	16.62	266	828	0.376	ng/ul	94
12) Phenanthrene	16.99	178	12451	0.373	ng/ul	100
13) Anthracene	17.08	178	10505	0.376	ng/ul	99
15) Fluoranthene	19.01	202	14197	0.367	ng/ul	100
17) Pyrene	19.38	202	14117	0.380	ng/ul	98
18) Benzo(a)anthracene	21.13	228	12293	0.389	ng/ul	99
19) Chrysene	21.18	228	14959	0.392	ng/ul	99
21) Benzo(b)fluoranthene	22.66	252	13887	0.393	ng/ul	100
22) Benzo(k)fluoranthene	22.70	252	14822	0.409	ng/ul	99
23) Benzo(a)pyrene	23.20	252	12309	0.423	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.42	276	16817	0.455	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.43	278	13811	0.483	ng/ul	96
26) Benzo(g,h,i)perylene	26.06	276	14853	0.413	ng/ul	99

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

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