

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM081720\
 Data File : BM027115.D
 Acq On : 17 Aug 2020 14:58
 Operator : JU/CG
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SICV82

Quant Time: Aug 17 15:56:41 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM081720.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 17 14:20:13 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1554	0.40	ng/ul	0.00
2) Naphthalene-d8	10.40	136	4852	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.27	164	2917	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	6611	0.40	ng/ul	0.00
16) Chrysene-d12	21.21	240	6276	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	6309	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	2816	0.34	ng/ul	0.00
14) Fluoranthene-d10	19.04	212	6448	0.34	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.45	128	5148	0.354	ng/ul	99
5) 2-Methylnaphthalene	12.07	142	3702	0.372	ng/ul	97
7) Acenaphthylene	13.99	152	4516	0.357	ng/ul	99
8) Acenaphthene	14.33	153	4048	0.368	ng/ul	99
9) Fluorene	15.32	166	4826	0.366	ng/ul	99
11) Pentachlorophenol	16.67	266	743	0.341	ng/ul	100
12) Phenanthrene	17.05	178	8203	0.367	ng/ul	99
13) Anthracene	17.15	178	7146	0.352	ng/ul	99
15) Fluoranthene	19.07	202	9610	0.358	ng/ul	98
17) Pyrene	19.44	202	9663	0.360	ng/ul	100
18) Benzo(a)anthracene	21.19	228	8747	0.351	ng/ul	99
19) Chrysene	21.24	228	10017	0.371	ng/ul	99
21) Benzo(b)fluoranthene	22.75	252	10089	0.373	ng/ul	100
22) Benzo(k)fluoranthene	22.79	252	10163	0.361	ng/ul	100
23) Benzo(a)pyrene	23.30	252	8917	0.364	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.59	276	12269	0.363	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.60	278	10087	0.367	ng/ul	99
26) Benzo(g,h,i)perylene	26.26	276	10426	0.357	ng/ul	98

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

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