

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM111020\
 Data File : BM027772.D
 Acq On : 10 Nov 2020 16:27
 Operator : JU/CG
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SICV82

Quant Time: Nov 10 16:58:53 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM111020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 10 15:45:47 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.39	152	538	0.40	ng/ul	0.00
2) Naphthalene-d8	11.22	136	1924	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.99	164	1224	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	2616	0.40	ng/ul	0.00
16) Chrysene-d12	21.81	240	2054	0.40	ng/ul	0.00
20) Perylene-d12	24.25	264	1631	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.79	152	1321	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	2763	0.36	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.27	128	2306	0.375	ng/ul	100
5) 2-Methylnaphthalene	12.86	142	1652	0.386	ng/ul	100
7) Acenaphthylene	14.72	152	2287	0.380	ng/ul	99
8) Acenaphthene	15.06	153	1727	0.374	ng/ul	99
9) Fluorene	16.03	166	2066	0.372	ng/ul	99
11) Pentachlorophenol	17.35	266	337	0.357	ng/ul	97
12) Phenanthrene	17.75	178	3181	0.374	ng/ul	99
13) Anthracene	17.84	178	3073	0.371	ng/ul	99
15) Fluoranthene	19.73	202	3578	0.367	ng/ul	99
17) Pyrene	20.08	202	3544	0.389	ng/ul	98
18) Benzo(a)anthracene	21.79	228	3116	0.380	ng/ul	98
19) Chrysene	21.85	228	3202	0.383	ng/ul	99
21) Benzo(b)fluoranthene	23.50	252	3005	0.403	ng/ul	97
22) Benzo(k)fluoranthene	23.55	252	2869	0.392	ng/ul	96
23) Benzo(a)pyrene	24.14	252	2459	0.375	ng/ul	94
24) Indeno(1,2,3-cd)pyrene	26.77	276	2806	0.375	ng/ul#	99
25) Dibenzo(a,h)anthracene	26.79	278	2334	0.374	ng/ul#	94
26) Benzo(g,h,i)perylene	27.55	276	2370	0.375	ng/ul	98

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

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