

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM020521\
 Data File : BM028668.D
 Acq On : 05 Feb 2021 17:35
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SICV82

Quant Time: Feb 05 18:08:21 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM020521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 05 17:33:51 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	471	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	1929	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.56	164	999	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.29	188	1907	0.40	ng/ul	0.00
16) Chrysene-d12	21.44	240	1705	0.40	ng/ul	0.00
20) Perylene-d12	23.68	264	1759	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.30	152	1139	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.31	212	2013	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.76	128	2020	0.421	ng/ul	99
5) 2-Methylnaphthalene	12.38	142	1323	0.411	ng/ul	98
7) Acenaphthylene	14.28	152	1656	0.406	ng/ul	100
8) Acenaphthene	14.62	153	1316	0.407	ng/ul	100
9) Fluorene	15.60	166	1441	0.402	ng/ul	98
11) Pentachlorophenol	16.95	266	197	0.395	ng/ul	95
12) Phenanthrene	17.33	178	2205	0.414	ng/ul	100
13) Anthracene	17.43	178	2078	0.409	ng/ul	100
15) Fluoranthene	19.34	202	2443	0.403	ng/ul	99
17) Pyrene	19.69	202	2521	0.407	ng/ul	99
18) Benzo(a)anthracene	21.42	228	2378	0.409	ng/ul	99
19) Chrysene	21.47	228	2383	0.412	ng/ul	99
21) Benzo(b)fluoranthene	23.00	252	2495	0.414	ng/ul	96
22) Benzo(k)fluoranthene	23.05	252	2448	0.407	ng/ul	96
23) Benzo(a)pyrene	23.58	252	2213	0.413	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.95	276	2865	0.410	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.96	278	2387	0.407	ng/ul	96
26) Benzo(g,h,i)perylene	26.65	276	2407	0.414	ng/ul	99

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

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