

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022420\
 Data File : BM025034.D
 Acq On : 24 Feb 2020 22:36
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SICV82

Quant Time: Feb 25 01:13:34 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM022420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 25 01:09:25 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	2110	0.40	ng/ul	0.00
2) Naphthalene-d8	10.47	136	8262	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.33	164	5757	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.07	188	13630	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	15317	0.40	ng/ul	0.00
20) Perylene-d12	23.45	264	16533	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.06	152	6353	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	19084	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.52	128	10224	0.407	ng/ul	93
5) 2-Methylnaphthalene	12.14	142	7480	0.412	ng/ul	100
7) Acenaphthylene	14.04	152	13147	0.396	ng/ul	97
8) Acenaphthene	14.39	153	8711	0.391	ng/ul	98
9) Fluorene	15.38	166	10912	0.390	ng/ul	99
11) Pentachlorophenol	16.72	266	1713	0.411	ng/ul	97
12) Phenanthrene	17.11	178	18668	0.389	ng/ul	98
13) Anthracene	17.20	178	18608	0.400	ng/ul	98
15) Fluoranthene	19.13	202	25558	0.367	ng/ul	98
17) Pyrene	19.49	202	25564	0.370	ng/ul	99
18) Benzo(a)anthracene	21.24	228	25971	0.393	ng/ul	98
19) Chrysene	21.29	228	26173	0.397	ng/ul	99
21) Benzo(b)fluoranthene	22.80	252	26666	0.411	ng/ul	96
22) Benzo(k)fluoranthene	22.84	252	25914	0.404	ng/ul	98
23) Benzo(a)pyrene	23.36	252	24218	0.409	ng/ul#	95
24) Indeno(1,2,3-cd)pyrene	25.65	276	30425	0.410	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.67	278	25350	0.407	ng/ul	97
26) Benzo(g,h,i)perylene	26.31	276	25520	0.409	ng/ul	97

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

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