

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM041520\
 Data File : BM025962.D
 Acq On : 15 Apr 2020 12:28
 Operator : MA/SJ
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SICV82

Quant Time: Apr 15 12:59:27 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM041520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 15 12:24:52 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	1623	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	6005	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	3182	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	7490	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	5574	0.40	ng/ul	0.00
20) Perylene-d12	23.28	264	4523	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.92	152	3781	0.38	ng/ul	0.00
14) Fluoranthene-d10	18.97	212	8484	0.39	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.37	128	6983	0.370	ng/ul	99
5) 2-Methylnaphthalene	11.99	142	4795	0.383	ng/ul	98
7) Acenaphthylene	13.91	152	6355	0.428	ng/ul	98
8) Acenaphthene	14.26	153	5073	0.404	ng/ul	99
9) Fluorene	15.25	166	5959	0.405	ng/ul	99
11) Pentachlorophenol	16.61	266	685	0.362	ng/ul	97
12) Phenanthrene	16.98	178	9698	0.380	ng/ul	99
13) Anthracene	17.08	178	8101	0.377	ng/ul	98
15) Fluoranthene	19.00	202	10960	0.369	ng/ul	98
17) Pyrene	19.37	202	10980	0.409	ng/ul	98
18) Benzo(a)anthracene	21.12	228	8116	0.385	ng/ul	99
19) Chrysene	21.17	228	10022	0.383	ng/ul	100
21) Benzo(b)fluoranthene	22.65	252	7910	0.383	ng/ul	100
22) Benzo(k)fluoranthene	22.69	252	9027	0.397	ng/ul	98
23) Benzo(a)pyrene	23.19	252	7269	0.420	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.41	276	8174	0.429	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.42	278	6717	0.458	ng/ul	96
26) Benzo(g,h,i)perylene	26.04	276	7042	0.403	ng/ul	97

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

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