

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030920\
 Data File : BM025415.D
 Acq On : 09 Mar 2020 13:34
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SICV79

Quant Time: Mar 09 14:43:00 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM030920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 09 13:32:56 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	2669	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	10344	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	6857	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	14899	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	11103	0.40	ng/ul	0.00
20) Perylene-d12	23.41	264	9976	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	6263	0.34	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	12957	0.32	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.48	128	10419	0.340	ng/ul	100
5) 2-Methylnaphthalene	12.10	142	7308	0.334	ng/ul	99
7) Acenaphthylene	14.01	152	9274	0.331	ng/ul	99
8) Acenaphthene	14.35	153	8128	0.331	ng/ul	99
9) Fluorene	15.34	166	9712	0.324	ng/ul	100
11) Pentachlorophenol	16.70	266	1052	0.301	ng/ul	97
12) Phenanthrene	17.08	178	15426	0.327	ng/ul	100
13) Anthracene	17.17	178	13048	0.318	ng/ul	100
15) Fluoranthene	19.09	202	16189	0.316	ng/ul	99
17) Pyrene	19.45	202	15750	0.316	ng/ul	99
18) Benzo(a)anthracene	21.21	228	12758	0.321	ng/ul	99
19) Chrysene	21.26	228	14866	0.333	ng/ul	99
21) Benzo(b)fluoranthene	22.75	252	13268	0.336	ng/ul	99
22) Benzo(k)fluoranthene	22.80	252	13627	0.329	ng/ul	98
23) Benzo(a)pyrene	23.31	252	10671	0.325	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.58	276	14563	0.327	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.59	278	12106	0.323	ng/ul	99
26) Benzo(g,h,i)perylene	26.24	276	12572	0.326	ng/ul	99

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

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