

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040521\
 Data File : BM029344.D
 Acq On : 05 Apr 2021 16:32
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SICV82

Quant Time: Apr 05 17:03:32 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM040521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 05 15:51:17 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	1944	0.40	ng/ul	0.00
2) Naphthalene-d8	10.48	136	6315	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.33	164	3160	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.07	188	6408	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	6590	0.40	ng/ul	0.00
20) Perylene-d12	23.46	264	6875	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.07	152	3534	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.10	212	6447	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.53	128	7156	0.418	ng/ul	98
5) 2-Methylnaphthalene	12.15	142	4510	0.407	ng/ul	98
7) Acenaphthylene	14.05	152	5430	0.405	ng/ul	99
8) Acenaphthene	14.39	153	4680	0.412	ng/ul	97
9) Fluorene	15.38	166	5152	0.410	ng/ul	99
11) Pentachlorophenol	16.73	266	547	0.377	ng/ul	93
12) Phenanthrene	17.11	178	7824	0.400	ng/ul	99
13) Anthracene	17.20	178	6848	0.399	ng/ul	99
15) Fluoranthene	19.13	202	8099	0.373	ng/ul	99
17) Pyrene	19.49	202	8351	0.393	ng/ul	99
18) Benzo(a)anthracene	21.23	228	7085	0.377	ng/ul	100
19) Chrysene	21.28	228	8028	0.389	ng/ul	99
21) Benzo(b)fluoranthene	22.79	252	8532	0.409	ng/ul	98
22) Benzo(k)fluoranthene	22.84	252	8722	0.408	ng/ul	97
23) Benzo(a)pyrene	23.36	252	7280	0.403	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.68	276	10133	0.405	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.70	278	8552	0.408	ng/ul	100
26) Benzo(g,h,i)perylene	26.36	276	8724	0.411	ng/ul	99

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

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