

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM080319\
 Data File : BM021785.D
 Acq On : 03 Aug 2019 21:08
 Operator : HP/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SICV25

Quant Time: Aug 03 23:08:02 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM080319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Aug 03 22:53:27 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	704	0.40	ng/ul	0.00
2) Naphthalene-d8	10.44	136	2848	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	1582	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.07	188	3589	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	3529	0.40	ng/ul	0.00
20) Perylene-d12	23.48	264	4142	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.06	152	1790	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	4698	0.39	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.50	128	3130	0.389	ng/ul	99
5) 2-Methylnaphthalene	12.13	142	2064	0.385	ng/ul	100
7) Acenaphthylene	14.03	152	2909	0.394	ng/ul	99
8) Acenaphthene	14.37	153	2375	0.393	ng/ul	97
9) Fluorene	15.37	166	2584	0.385	ng/ul	99
11) Pentachlorophenol	16.75	266	223	0.337	ng/ul	96
12) Phenanthrene	17.11	178	4111	0.394	ng/ul	99
13) Anthracene	17.21	178	3329	0.374	ng/ul	98
15) Fluoranthene	19.12	202	5402	0.386	ng/ul	98
17) Pyrene	19.49	202	5832	0.402	ng/ul	98
18) Benzo(a)anthracene	21.24	228	4708	0.373	ng/ul	99
19) Chrysene	21.29	228	6529	0.402	ng/ul	99
21) Benzo(b)fluoranthene	22.81	252	5426	0.388	ng/ul	99
22) Benzo(k)fluoranthene	22.85	252	6290	0.398	ng/ul	99
23) Benzo(a)pyrene	23.39	252	5542	0.390	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.70	276	6610	0.390	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.73	278	5242	0.382	ng/ul	97
26) Benzo(g,h,i)perylene	26.38	276	6057	0.397	ng/ul	97

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

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