

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE102518\
 Data File : BE098056.D
 Acq On : 25 Oct 2018 15:28
 Operator : SJ/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SICV67

Quant Time: Oct 25 16:41:05 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE102518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 25 14:48:33 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.88	152	889	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	4614	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	3290	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.29	188	8188	0.40	ng/ul	0.00
16) Chrysene-d12	21.47	240	10989	0.40	ng/ul	0.00
20) Perylene-d12	24.02	264	9582	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	3264	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	10399	0.39	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.73	128	4304	0.390	ng/ul	98
5) 2-Methylnaphthalene	12.34	142	3275	0.392	ng/ul	100
7) Acenaphthylene	14.25	152	5037	0.388	ng/ul	99
8) Acenaphthene	14.60	153	3692	0.381	ng/ul	99
9) Fluorene	15.59	166	4565	0.391	ng/ul	98
11) Pentachlorophenol	16.95	266	155	0.362	ng/ul	94
12) Phenanthrene	17.33	178	7421	0.393	ng/ul	98
13) Anthracene	17.42	178	7170	0.384	ng/ul	100
15) Fluoranthene	19.35	202	9348	0.388	ng/ul	99
17) Pyrene	19.71	202	9545	0.372	ng/ul	98
18) Benzo(a)anthracene	21.45	228	13604	0.421	ng/ul	99
19) Chrysene	21.51	228	10364	0.415	ng/ul	99
21) Benzo(b)fluoranthene	23.24	252	10549	0.397	ng/ul	99
22) Benzo(k)fluoranthene	23.29	252	10261	0.395	ng/ul	98
23) Benzo(a)pyrene	23.91	252	10015	0.392	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	26.71	276	11699	0.400	ng/ul#	98
25) Dibenzo(a,h)anthracene	26.74	278	9717	0.398	ng/ul	97
26) Benzo(g,h,i)perylene	27.54	276	9758	0.396	ng/ul	97

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

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