

Data Path : Z:\HPCHEM1\BNA_E\Data\BE111417\
 Data File : BE094808.D
 Acq On : 14 Nov 2017 15:03
 Operator : SJ/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SICV90

Quant Time: Nov 14 16:35:04 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE111417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 14 16:33:16 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.93	152	708	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	4122	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	2561	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.27	188	6826	0.40	ng/ul	0.00
16) Chrysene-d12	21.44	240	8344	0.40	ng/ul	0.00
20) Perylene-d12	23.98	264	7847	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.31	152	2658	0.40	ng/ul	0.01
14) Fluoranthene-d10	19.29	212	9458	0.40	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.76	128	3921	0.399	ng/ul#	94
5) 2-Methylnaphthalene	12.39	142	2721	0.398	ng/ul	99
7) Acenaphthylene	14.25	152	4607	0.393	ng/ul	98
8) Acenaphthene	14.59	153	3306	0.401	ng/ul	97
9) Fluorene	15.58	166	4039	0.403	ng/ul	94
11) Pentachlorophenol	17.05	266	131	0.364	ng/ul	94
12) Phenanthrene	17.31	178	7302	0.398	ng/ul#	90
13) Anthracene	17.41	178	7382	0.395	ng/ul#	93
15) Fluoranthene	19.33	202	9708	0.404	ng/ul	84
17) Pyrene	19.69	202	10381	0.398	ng/ul#	82
18) Benzo(a)anthracene	21.42	228	10006	0.394	ng/ul#	85
19) Chrysene	21.48	228	10194	0.403	ng/ul	99
21) Benzo(b)fluoranthene	23.19	252	9383	0.401	ng/ul	90
22) Benzo(k)fluoranthene	23.25	252	9702	0.390	ng/ul	91
23) Benzo(a)pyrene	23.87	252	9175	0.395	ng/ul#	86
24) Indeno(1,2,3-cd)pyrene	26.70	276	9979	0.394	ng/ul#	81
25) Dibenzo(a,h)anthracene	26.69	278	8393	0.393	ng/ul	94
26) Benzo(g,h,i)perylene	27.55	276	8276	0.390	ng/ul	96

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

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