

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE092017\
 Data File : BE094030.D
 Acq On : 20 Sep 2017 12:40
 Operator : SJ/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SICV08

Quant Time: Sep 20 13:29:29 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 20 13:07:25 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1308	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	6048	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	3497	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.27	188	9851	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	12824	0.40	ng/ul	0.00
20) Perylene-d12	23.94	264	11400	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.32	152	3795	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	12833	0.39	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.75	128	5977	0.403	ng/ul#	94
5) 2-Methylnaphthalene	12.39	142	3866	0.402	ng/ul	100
7) Acenaphthylene	14.27	152	5429	0.398	ng/ul	98
8) Acenaphthene	14.60	153	4501	0.401	ng/ul	97
9) Fluorene	15.59	166	5317	0.405	ng/ul	91
11) Pentachlorophenol	16.96	266	376	0.325	ng/ul	92
12) Phenanthrene	17.31	178	10578	0.411	ng/ul#	88
13) Anthracene	17.41	178	9682	0.387	ng/ul#	93
15) Fluoranthene	19.32	202	13287	0.393	ng/ul	86
17) Pyrene	19.69	202	14417	0.404	ng/ul#	83
18) Benzo(a)anthracene	21.42	228	13383	0.393	ng/ul#	87
19) Chrysene	21.47	228	16096	0.410	ng/ul	98
21) Benzo(b)fluoranthene	23.17	252	12264	0.396	ng/ul	86
22) Benzo(k)fluoranthene	23.22	252	14560	0.407	ng/ul#	89
23) Benzo(a)pyrene	23.83	252	12787	0.397	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	26.61	276	13121	0.399	ng/ul#	84
25) Dibenzo(a,h)anthracene	26.63	278	11049	0.400	ng/ul	92
26) Benzo(g,h,i)perylene	27.42	276	11418	0.402	ng/ul	99

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

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