

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE083118\
 Data File : BE097407.D
 Acq On : 31 Aug 2018 17:47
 Operator : SJ/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.423

Quant Time: Aug 31 18:25:33 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE082918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 31 18:03:11 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	797	0.40	ng/ul	0.00
2) Naphthalene-d8	10.12	136	4011	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.01	164	2387	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.76	188	6689	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	8162	0.40	ng/ul	0.00
20) Perylene-d12	23.20	264	8057	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	2603	0.42	ng/ul	0.00
14) Fluoranthene-d10	18.80	212	8839	0.40	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.17	128	3942	0.403	ng/ul#	94
5) 2-Methylnaphthalene	11.80	142	2599	0.420	ng/ul	98
7) Acenaphthylene	13.72	152	3664	0.421	ng/ul	99
8) Acenaphthene	14.07	153	3026	0.405	ng/ul	94
9) Fluorene	15.07	166	3557	0.415	ng/ul	93
11) Pentachlorophenol	16.44	266	302	0.458	ng/ul	98
12) Phenanthrene	16.80	178	7021	0.408	ng/ul#	89
13) Anthracene	16.90	178	6155	0.425	ng/ul#	92
15) Fluoranthene	18.83	202	8789	0.397	ng/ul	84
17) Pyrene	19.20	202	8828	0.402	ng/ul#	81
18) Benzo(a)anthracene	20.95	228	9025	0.430	ng/ul#	85
19) Chrysene	21.01	228	9799	0.406	ng/ul	97
21) Benzo(b)fluoranthene	22.53	252	9091	0.421	ng/ul	84
22) Benzo(k)fluoranthene	22.57	252	9269	0.385	ng/ul#	89
23) Benzo(a)pyrene	23.10	252	8505	0.426	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	25.49	276	10903	0.421	ng/ul#	77
25) Dibenzo(a,h)anthracene	25.51	278	9148	0.417	ng/ul#	91
26) Benzo(g,h,i)perylene	26.19	276	9234	0.419	ng/ul#	92

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

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