

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE090718\
 Data File : BE097416.D
 Acq On : 7 Sep 2018 9:23
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.426

Quant Time: Sep 07 23:36:16 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 : Thu Sep 06 01:29:05 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	816	0.40	ng/ul	0.00
2) Naphthalene-d8	10.12	136	3989	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.01	164	2356	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.76	188	6990	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	7783	0.40	ng/ul	0.00
20) Perylene-d12	23.20	264	7022	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	11.72	152	2553	0.41	ng/ul	0.00
14) Fluoranthene-d10	18.80	212	8733	0.38	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.17	128	3906	0.402	ng/ul#	94
5) 2-Methylnaphthalene	11.80	142	2532	0.412	ng/ul	97
7) Acenaphthylene	13.72	152	3508	0.409	ng/ul	98
8) Acenaphthene	14.07	153	2980	0.404	ng/ul	95
9) Fluorene	15.07	166	3491	0.413	ng/ul	94
11) Pentachlorophenol	16.44	266	252	0.366	ng/ul	97
12) Phenanthrene	16.80	178	7259	0.403	ng/ul#	90
13) Anthracene	16.90	178	6153	0.407	ng/ul#	91
15) Fluoranthene	18.83	202	8786	0.380	ng/ul	84
17) Pyrene	19.20	202	8510	0.406	ng/ul#	81
18) Benzo(a)anthracene	20.95	228	7961	0.398	ng/ul#	84
19) Chrysene	21.01	228	9426	0.409	ng/ul	98
21) Benzo(b)fluoranthene	22.53	252	7772	0.413	ng/ul	83
22) Benzo(k)fluoranthene	22.57	252	8483	0.405	ng/ul#	86
23) Benzo(a)pyrene	23.10	252	7107	0.409	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	25.49	276	9227	0.409	ng/ul#	75
25) Dibenzo(a,h)anthracene	25.51	278	7674	0.402	ng/ul#	87
26) Benzo(g,h,i)perylene	26.19	276	8023	0.417	ng/ul#	92

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

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