

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE090518\
 Data File : BE097414.D
 Acq On : 5 Sep 2018 20:20
 Operator : SJ/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.425

Quant Time: Sep 06 01:30:16 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	914	0.40	ng/ul	0.00
2) Naphthalene-d8	10.12	136	4484	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.01	164	2558	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.76	188	6906	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	7729	0.40	ng/ul	0.00
20) Perylene-d12	23.20	264	7276	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	2854	0.41	ng/ul	0.00
14) Fluoranthene-d10	18.80	212	8711	0.38	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.17	128	4373	0.400	ng/ul#	93
5) 2-Methylnaphthalene	11.79	142	2845	0.412	ng/ul	97
7) Acenaphthylene	13.72	152	3835	0.412	ng/ul	99
8) Acenaphthene	14.07	153	3139	0.392	ng/ul	94
9) Fluorene	15.07	166	3672	0.400	ng/ul	91
11) Pentachlorophenol	16.44	266	255	0.375	ng/ul	97
12) Phenanthrene	16.80	178	7165	0.403	ng/ul#	90
13) Anthracene	16.90	178	6190	0.414	ng/ul#	92
15) Fluoranthene	18.83	202	8684	0.380	ng/ul	84
17) Pyrene	19.20	202	8465	0.407	ng/ul#	79
18) Benzo(a)anthracene	20.95	228	8090	0.407	ng/ul#	86
19) Chrysene	21.00	228	9211	0.403	ng/ul	95
21) Benzo(b)fluoranthene	22.53	252	7816	0.401	ng/ul	84
22) Benzo(k)fluoranthene	22.57	252	8725	0.402	ng/ul#	86
23) Benzo(a)pyrene	23.10	252	7432	0.412	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	25.49	276	9513	0.407	ng/ul#	76
25) Dibenzo(a,h)anthracene	25.51	278	7987	0.403	ng/ul#	88
26) Benzo(g,h,i)perylene	26.19	276	8135	0.408	ng/ul#	92

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

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