

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE101518\
 Data File : BE097903.D
 Acq On : 15 Oct 2018 17:18
 Operator : SJ/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.457

Quant Time: Oct 15 17:48:18 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE101118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 15 16:15:33 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1905	0.40	ng/ul	0.00
2) Naphthalene-d8	10.72	136	9922	0.40	ng/ul	0.01
6) Acenaphthene-d10	14.57	164	6298	0.40	ng/ul	0.01
10) Phenanthrene-d10	17.32	188	14696	0.40	ng/ul	0.00
16) Chrysene-d12	21.50	240	14897	0.40	ng/ul	0.00
20) Perylene-d12	24.07	264	15348	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.31	152	7242	0.43	ng/ul	0.01
14) Fluoranthene-d10	19.35	212	20232	0.43	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.76	128	10469	0.434	ng/ul	98
5) 2-Methylnaphthalene	12.38	142	7512	0.430	ng/ul	100
7) Acenaphthylene	14.29	152	11367	0.477	ng/ul	97
8) Acenaphthene	14.63	153	8185	0.443	ng/ul	97
9) Fluorene	15.62	166	9532	0.395	ng/ul	98
11) Pentachlorophenol	16.99	266	1120	0.541	ng/ul#	84
12) Phenanthrene	17.36	178	15224	0.438	ng/ul	96
13) Anthracene	17.45	178	15346	0.476	ng/ul	98
15) Fluoranthene	19.38	202	19014	0.426	ng/ul	98
17) Pyrene	19.74	202	23500	0.530	ng/ul	96
18) Benzo(a)anthracene	21.49	228	19959	0.443	ng/ul#	91
19) Chrysene	21.54	228	19392	0.479	ng/ul#	88
21) Benzo(b)fluoranthene	23.29	252	19220	0.469	ng/ul#	54
22) Benzo(k)fluoranthene	23.34	252	19454	0.443	ng/ul#	45
23) Benzo(a)pyrene	23.96	252	18723	0.465	ng/ul#	31
24) Indeno(1,2,3-cd)pyrene	26.79	276	20979	0.459	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.82	278	17619	0.466	ng/ul#	53
26) Benzo(g,h,i)perylene	27.63	276	16515	0.415	ng/ul#	66

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

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