

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE071118\
 Data File : BE096893.D
 Acq On : 11 Jul 2018 17:07
 Operator : SJ/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.456

Quant Time: Jul 11 18:10:55 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE070618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 10 02:37:30 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	2571	0.40	ng/ul	0.00
2) Naphthalene-d8	10.14	136	14234	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.03	164	7962	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.78	188	18768	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	16339	0.40	ng/ul	0.00
20) Perylene-d12	23.22	264	17113	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.74	152	7977	0.34	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	15210	0.33	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.19	128	11533	0.335	ng/ul#	91
5) 2-Methylnaphthalene	11.81	142	8059	0.344	ng/ul	98
7) Acenaphthylene	13.74	152	11130	0.341	ng/ul	96
8) Acenaphthene	14.09	153	8655	0.321	ng/ul	95
9) Fluorene	15.08	166	8728	0.314	ng/ul	93
11) Pentachlorophenol	16.44	266	467	0.380	ng/ul	92
12) Phenanthrene	16.82	178	15490	0.320	ng/ul#	89
13) Anthracene	16.91	178	13713	0.334	ng/ul#	92
15) Fluoranthene	18.84	202	16704	0.324	ng/ul	85
17) Pyrene	19.21	202	16451	0.307	ng/ul#	81
18) Benzo(a)anthracene	20.96	228	14062	0.365	ng/ul#	85
19) Chrysene	21.02	228	15798	0.316	ng/ul	97
21) Benzo(b)fluoranthene	22.54	252	14532	0.311	ng/ul	82
22) Benzo(k)fluoranthene	22.58	252	14711	0.288	ng/ul#	85
23) Benzo(a)pyrene	23.11	252	13099	0.332	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	25.51	276	14616	0.356	ng/ul#	80
25) Dibenzo(a,h)anthracene	25.53	278	12115	0.372	ng/ul#	88
26) Benzo(g,h,i)perylene	26.21	276	12871	0.323	ng/ul#	93

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

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