

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE071718\
 Data File : BE096985.D
 Acq On : 18 Jul 2018 8:19
 Operator : SJ/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.470

Quant Time: Jul 18 09:04:59 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE071718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 18 07:15:45 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	1441	0.40	ng/ul	0.00
2) Naphthalene-d8	10.14	136	6964	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.03	164	4086	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.78	188	11775	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	12902	0.40	ng/ul	0.00
20) Perylene-d12	23.22	264	13192	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.74	152	4619	0.42	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	13879	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.18	128	6780	0.402	ng/ul#	93
5) 2-Methylnaphthalene	11.81	142	4651	0.418	ng/ul	97
7) Acenaphthylene	13.74	152	7440	0.418	ng/ul	97
8) Acenaphthene	14.09	153	5534	0.400	ng/ul	94
9) Fluorene	15.08	166	6142	0.418	ng/ul	95
11) Pentachlorophenol	16.43	266	667	0.501	ng/ul	96
12) Phenanthrene	16.81	178	12120	0.392	ng/ul#	88
13) Anthracene	16.90	178	11204	0.415	ng/ul#	91
15) Fluoranthene	18.84	202	15213	0.373	ng/ul	84
17) Pyrene	19.21	202	15886	0.426	ng/ul#	82
18) Benzo(a)anthracene	20.96	228	14378	0.421	ng/ul#	85
19) Chrysene	21.01	228	15143	0.391	ng/ul	97
21) Benzo(b)fluoranthene	22.53	252	15076	0.403	ng/ul	84
22) Benzo(k)fluoranthene	22.58	252	14399	0.372	ng/ul#	86
23) Benzo(a)pyrene	23.11	252	13615	0.414	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	25.50	276	15738	0.427	ng/ul#	80
25) Dibenzo(a,h)anthracene	25.53	278	13025	0.423	ng/ul#	88
26) Benzo(g,h,i)perylene	26.20	276	13538	0.412	ng/ul	94

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

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