

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE071118\
 Data File : BE096897.D
 Acq On : 12 Jul 2018 13:35
 Operator : SJ/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.458

Quant Time: Jul 12 15:08:06 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	2611	0.40	ng/ul	0.00
2) Naphthalene-d8	10.14	136	14475	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.03	164	8261	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.78	188	19708	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	17033	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	17115	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.74	152	8166	0.34	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	15702	0.32	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.19	128	11819	0.338	ng/ul#	92
5) 2-Methylnaphthalene	11.81	142	8321	0.349	ng/ul	97
7) Acenaphthylene	13.74	152	11506	0.340	ng/ul	96
8) Acenaphthene	14.09	153	8966	0.321	ng/ul	96
9) Fluorene	15.08	166	9041	0.313	ng/ul	91
11) Pentachlorophenol	16.43	266	452	0.350	ng/ul	98
12) Phenanthrene	16.82	178	16342	0.322	ng/ul#	89
13) Anthracene	16.90	178	14205	0.329	ng/ul#	91
15) Fluoranthene	18.84	202	17438	0.322	ng/ul	84
17) Pyrene	19.21	202	16997	0.304	ng/ul#	78
18) Benzo(a)anthracene	20.96	228	13947	0.347	ng/ul#	85
19) Chrysene	21.02	228	16489	0.317	ng/ul	98
21) Benzo(b)fluoranthene	22.54	252	14449	0.309	ng/ul	82
22) Benzo(k)fluoranthene	22.58	252	14691	0.288	ng/ul#	85
23) Benzo(a)pyrene	23.11	252	12672	0.321	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	25.51	276	13986	0.341	ng/ul#	80
25) Dibenzo(a,h)anthracene	25.53	278	11575	0.355	ng/ul#	88
26) Benzo(g,h,i)perylene	26.21	276	12427	0.312	ng/ul#	93

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

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