

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE071718\
 Data File : BE096976.D
 Acq On : 18 Jul 2018 2:42
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.469

Quant Time: Jul 18 05:38:21 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 03:40:41 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	1483	0.40	ng/ul	0.00
2) Naphthalene-d8	10.14	136	6963	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	3830	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.78	188	11025	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	12525	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	12540	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.73	152	4550	0.41	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	13283	0.40	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
3) Naphthalene	10.19	128	6788	0.402	ng/ul#	93
5) 2-Methylnaphthalene	11.81	142	4562	0.410	ng/ul	97
7) Acenaphthylene	13.74	152	6799	0.407	ng/ul	98
8) Acenaphthene	14.09	153	5131	0.396	ng/ul	94
9) Fluorene	15.08	166	5626	0.409	ng/ul	95
11) Pentachlorophenol	16.43	266	567	0.455	ng/ul	94
12) Phenanthrene	16.81	178	11372	0.393	ng/ul#	89
13) Anthracene	16.90	178	10293	0.407	ng/ul#	92
15) Fluoranthene	18.84	202	14479	0.379	ng/ul	84
17) Pyrene	19.21	202	15036	0.415	ng/ul#	81
18) Benzo(a)anthracene	20.96	228	13602	0.410	ng/ul#	85
19) Chrysene	21.01	228	14484	0.385	ng/ul	96
21) Benzo(b)fluoranthene	22.54	252	13071	0.367	ng/ul	83
22) Benzo(k)fluoranthene	22.58	252	14191	0.386	ng/ul#	85
23) Benzo(a)pyrene	23.11	252	12266	0.393	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	25.50	276	14097	0.403	ng/ul#	81
25) Dibenzo(a,h)anthracene	25.52	278	11781	0.403	ng/ul#	88
26) Benzo(g,h,i)perylene	26.20	276	12305	0.394	ng/ul	95

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

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