

Data Path : Z:\HPCHEM1\BNA E\DATA\BE032818\
 Data File : BE095896.D
 Acq On : 29 Mar 2018 2:40
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.425

Quant Time: Mar 29 04:45:21 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE031318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 29 04:43:20 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.42	152	1597	0.40	ng/ul	0.00
2) Naphthalene-d8	11.24	136	4278	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.99	164	2636	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.70	188	6717	0.40	ng/ul	0.00
16) Chrysene-d12	21.79	240	7352	0.40	ng/ul	0.00
20) Perylene-d12	24.41	264	7129	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.79	152	3049	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.68	212	10039	0.47	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.29	128	4369	0.366	ng/ul#	89
5) 2-Methylnaphthalene	12.87	142	3082	0.371	ng/ul	100
7) Acenaphthylene	14.72	152	4984	0.374	ng/ul	97
8) Acenaphthene	15.06	153	3594	0.359	ng/ul	97
9) Fluorene	16.02	166	4138	0.380	ng/ul#	90
11) Pentachlorophenol	17.35	266	344	0.364	ng/ul	95
12) Phenanthrene	17.74	178	7256	0.361	ng/ul#	88
13) Anthracene	17.83	178	7132	0.379	ng/ul#	92
15) Fluoranthene	19.71	202	9644	0.396	ng/ul	84
17) Pyrene	20.06	202	10357	0.371	ng/ul#	79
18) Benzo(a)anthracene	21.77	228	9165	0.386	ng/ul#	85
19) Chrysene	21.83	228	8682	0.359	ng/ul	97
21) Benzo(b)fluoranthene	23.59	252	8457	0.342	ng/ul	87
22) Benzo(k)fluoranthene	23.65	252	8377	0.357	ng/ul#	88
23) Benzo(a)pyrene	24.29	252	8115	0.360	ng/ul#	81
24) Indeno(1,2,3-cd)pyrene	27.21	276	9312	0.372	ng/ul#	77
25) Dibenzo(a,h)anthracene	27.23	278	7535	0.377	ng/ul#	90
26) Benzo(g,h,i)perylene	28.09	276	7763	0.361	ng/ul	95

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

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