

Data Path : Z:\HPCHEM1\BNA E\DATA\BE012718\
 Data File : BE095162.D
 Acq On : 27 Jan 2018 12:24
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.413

Quant Time: Jan 28 01:33:15 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jan 27 03:18:43 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	2875	0.40	ng/ul	0.00
2) Naphthalene-d8	11.27	136	8548	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	5183	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.73	188	11909	0.40	ng/ul	0.02
16) Chrysene-d12	21.81	240	12937	0.40	ng/ul	0.00
20) Perylene-d12	24.45	264	12223	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.82	152	5873	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	17828	0.39	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	11.32	128	9539	0.399	ng/ul#	89
5) 2-Methylnaphthalene	12.89	142	6718	0.420	ng/ul	100
7) Acenaphthylene	14.75	152	10659	0.427	ng/ul	97
8) Acenaphthene	15.08	153	7695	0.398	ng/ul	97
9) Fluorene	16.03	166	8195	0.411	ng/ul#	69
11) Pentachlorophenol	17.36	266	1156	0.501	ng/ul	95
12) Phenanthrene	17.75	178	15861	0.396	ng/ul#	86
13) Anthracene	17.85	178	15911	0.434	ng/ul#	91
15) Fluoranthene	19.73	202	19686	0.380	ng/ul	84
17) Pyrene	20.08	202	19686	0.453	ng/ul#	81
18) Benzo(a)anthracene	21.79	228	18836	0.466	ng/ul#	86
19) Chrysene	21.84	228	16939	0.403	ng/ul	96
21) Benzo(b)fluoranthene	23.63	252	16754	0.389	ng/ul	91
22) Benzo(k)fluoranthene	23.68	252	15273	0.387	ng/ul	92
23) Benzo(a)pyrene	24.32	252	15452	0.401	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	27.25	276	17484	0.414	ng/ul#	78
25) Dibenzo(a,h)anthracene	27.28	278	14243	0.412	ng/ul#	93
26) Benzo(g,h,i)perylene	28.13	276	14478	0.405	ng/ul	96

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

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