

Data Path : U:\HPCHEM1\BNA E\DATA\BE011918\
 Data File : BE095078.D
 Acq On : 19 Jan 2018 16:26
 Operator : SJ/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SICV07

Quant Time: Jan 19 18:32:40 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE011918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 19 18:31:16 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.46	152	4603	0.40	ng/ul	0.00
2) Naphthalene-d8	11.29	136	14775	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.05	164	8197	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.75	188	18582	0.40	ng/ul	0.00
16) Chrysene-d12	21.84	240	18562	0.40	ng/ul	0.00
20) Perylene-d12	24.50	264	14798	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.84	152	9732	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.72	212	23262	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.34	128	15475	0.395	ng/ul#	89
5) 2-Methylnaphthalene	12.91	142	10858	0.398	ng/ul	100
7) Acenaphthylene	14.77	152	14282	0.389	ng/ul	96
8) Acenaphthene	15.11	153	11210	0.392	ng/ul	94
9) Fluorene	16.07	166	11641	0.382	ng/ul	90
11) Pentachlorophenol	17.41	266	365	0.332	ng/ul	97
12) Phenanthrene	17.78	178	19472	0.395	ng/ul#	81
13) Anthracene	17.87	178	17744	0.412	ng/ul#	87
15) Fluoranthene	19.76	202	25488	0.404	ng/ul	83
17) Pyrene	20.11	202	25434	0.377	ng/ul#	81
18) Benzo(a)anthracene	21.82	228	20298	0.395	ng/ul#	86
19) Chrysene	21.88	228	23123	0.404	ng/ul	96
21) Benzo(b)fluoranthene	23.67	252	19619	0.394	ng/ul	85
22) Benzo(k)fluoranthene	23.72	252	19386	0.422	ng/ul#	87
23) Benzo(a)pyrene	24.37	252	17399	0.409	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	27.33	276	16368	0.433	ng/ul#	81
25) Dibenzo(a,h)anthracene	27.35	278	12521	0.464	ng/ul#	89
26) Benzo(g,h,i)perylene	28.21	276	15384	0.410	ng/ul	94

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

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