

Data Path : Z:\HPCHEM1\BNA E\DATA\BE032818\
 Data File : BE095898.D
 Acq On : 29 Mar 2018 3:52
 Operator : SJ/JU
 Sample : J1951-04
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :

Quant Time: Mar 29 07:19:40 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 07:00:13 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.41	152	1899	0.40	ng/ul	0.00
2) Naphthalene-d8	11.24	136	5485	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.99	164	3612	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.70	188	8457m	0.40	ng/ul	0.00
16) Chrysene-d12	21.79	240	7971m	0.40	ng/ul	0.00
20) Perylene-d12	24.42	264	9104m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.79	152	2654	0.29	ng/ul	0.00
14) Fluoranthene-d10	19.68	212	7723m	0.29	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.29	128	3016	0.197	ng/ul#	91
5) 2-Methylnaphthalene	12.86	142	1609	0.151	ng/ul	99
7) Acenaphthylene	14.72	152	2076	0.114	ng/ul#	93
8) Acenaphthene	15.06	153	11990	0.873	ng/ul	96
9) Fluorene	16.02	166	11524	0.773	ng/ul#	90
12) Phenanthrene	17.75	178	274683m	10.845	ng/ul	
13) Anthracene	17.83	178	54939m	2.320	ng/ul	
15) Fluoranthene	19.71	202	628317m	20.480	ng/ul	
17) Pyrene	20.07	202	576307	19.041	ng/ul#	77
18) Benzo(a)anthracene	21.77	228	352684m	13.713	ng/ul	
19) Chrysene	21.83	228	263982	10.081	ng/ul	98
21) Benzo(b)fluoranthene	23.61	252	344366m	10.920	ng/ul	
22) Benzo(k)fluoranthene	23.65	252	129560m	4.322	ng/ul	
23) Benzo(a)pyrene	24.30	252	249017m	8.649	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.22	276	171087m	5.359	ng/ul	
25) Dibenzo(a,h)anthracene	27.23	278	43299m	1.696	ng/ul	
26) Benzo(a,h,i)perylene	28.10	276	159661m	5.822	ng/ul	

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

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