

Data Path : Z:\HPCHEM1\BNA E\DATA\BE042618\
 Data File : BE096189.D
 Acq On : 26 Apr 2018 21:08
 Operator : SJ/JU
 Sample : J2449-05
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :

Quant Time: Apr 27 05:55:40 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE042418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 27 04:45:57 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.36	152	681	0.40	ng/ul	0.00
2) Naphthalene-d8	11.18	136	2886	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.95	164	1745	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.66	188	3919m	0.40	ng/ul	0.00
16) Chrysene-d12	21.75	240	3528m	0.40	ng/ul	0.00
20) Perylene-d12	24.36	264	4420m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.74	152	1701	0.32	ng/ul	0.00
14) Fluoranthene-d10	19.64	212	4526m	0.25	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.23	128	11107	1.382	ng/ul#	88
5) 2-Methylnaphthalene	12.81	142	7239	1.345	ng/ul	100
7) Acenaphthylene	14.68	152	17707	1.968	ng/ul	97
8) Acenaphthene	15.01	153	5912	0.912	ng/ul	92
9) Fluorene	15.98	166	8937	1.041	ng/ul	92
12) Phenanthrene	17.70	178	227462m	19.714	ng/ul	
13) Anthracene	17.79	178	38543m	3.330	ng/ul	
15) Fluoranthene	19.67	202	581185m	32.953	ng/ul	
17) Pyrene	20.03	202	413365	58.404	ng/ul#	76
18) Benzo(a)anthracene	21.74	228	248677m	23.282	ng/ul	
19) Chrysene	21.79	228	250331	25.505	ng/ul	97
21) Benzo(b)fluoranthene	23.57	252	330911m	19.772	ng/ul	
22) Benzo(k)fluoranthene	23.61	252	124886m	7.806	ng/ul	
23) Benzo(a)pyrene	24.25	252	233284m	16.044	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.16	276	155098m	7.917	ng/ul	
25) Dibenzo(a,h)anthracene	27.16	278	42250m	2.576	ng/ul	
26) Benzo(a,h,i)perylene	28.03	276	120741m	7.080	ng/ul	

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

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