

Data Path : Z:\HPCHEM1\BNA E\DATA\BE042618\
 Data File : BE096190.D
 Acq On : 26 Apr 2018 21:43
 Operator : SJ/JU
 Sample : J2449-03
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :

Quant Time: Apr 27 06:01:05 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	770	0.40	ng/ul	0.00
2) Naphthalene-d8	11.18	136	3185	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.95	164	1886	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.66	188	4200m	0.40	ng/ul	0.00
16) Chrysene-d12	21.75	240	5274m	0.40	ng/ul	0.00
20) Perylene-d12	24.36	264	3944m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.74	152	2249	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.64	212	5884m	0.30	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	11.23	128	10156	1.145	ng/ul#	89
5) 2-Methylnaphthalene	12.81	142	5689	0.957	ng/ul	100
7) Acenaphthylene	14.68	152	4938	0.508	ng/ul	99
8) Acenaphthene	15.01	153	4057	0.579	ng/ul	91
9) Fluorene	15.98	166	7415	0.799	ng/ul	92
12) Phenanthrene	17.70	178	181094m	14.645	ng/ul	
13) Anthracene	17.79	178	32616m	2.629	ng/ul	
15) Fluoranthene	19.67	202	300467m	15.897	ng/ul	
17) Pyrene	20.03	202	211854	20.023	ng/ul#	77
18) Benzo(a)anthracene	21.73	228	97982m	6.137	ng/ul	
19) Chrysene	21.79	228	99615	6.789	ng/ul	97
21) Benzo(b)fluoranthene	23.55	252	123205m	8.250	ng/ul	
22) Benzo(k)fluoranthene	23.60	252	41664m	2.918	ng/ul	
23) Benzo(a)pyrene	24.25	252	89093	6.867	ng/ul#	77
24) Indeno(1,2,3-cd)pyrene	27.14	276	57874	3.311	ng/ul#	73
25) Dibenzo(a,h)anthracene	27.15	278	14623m	0.999	ng/ul	
26) Benzo(a,h,i)perylene	28.01	276	50538	3.321	ng/ul#	92

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

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