

Data Path : Z:\HPCHEM1\BNA E\DATA\BE042618\
 Data File : BE096191.D
 Acq On : 26 Apr 2018 22:19
 Operator : SJ/JU
 Sample : J2449-04
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AD9

Manual Integrations
 APPROVED

Sohil
 4/27/2018 7:29:34 PM

Quant Time: Apr 27 06:41:28 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	695	0.40	ng/ul	0.00
2) Naphthalene-d8	11.18	136	2831	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.95	164	1634	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.66	188	3772m	0.40	ng/ul	0.00
16) Chrysene-d12	21.75	240	4833m	0.40	ng/ul	0.00
20) Perylene-d12	24.36	264	3667	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.74	152	1134	0.22	ng/ul	0.00
14) Fluoranthene-d10	19.64	212	3047m	0.18	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.23	128	699	0.089	ng/ul#	76
5) 2-Methylnaphthalene	12.81	142	393	0.074	ng/ul	97
7) Acenaphthylene	14.68	152	337	0.040	ng/ul#	60
8) Acenaphthene	15.01	153	504	0.083	ng/ul#	93
9) Fluorene	15.98	166	655	0.082	ng/ul#	88
12) Phenanthrene	17.70	178	12129m	1.092	ng/ul	
13) Anthracene	17.79	178	2205m	0.198	ng/ul	
15) Fluoranthene	19.67	202	22139m	1.304	ng/ul	
17) Pyrene	20.02	202	12121	1.250	ng/ul#	79
18) Benzo(a)anthracene	21.73	228	9757	0.667	ng/ul#	86
19) Chrysene	21.79	228	8478	0.631	ng/ul	94
21) Benzo(b)fluoranthene	23.55	252	11369m	0.819	ng/ul	
22) Benzo(k)fluoranthene	23.60	252	3815m	0.287	ng/ul	
23) Benzo(a)pyrene	24.24	252	8103	0.672	ng/ul#	83
24) Indeno(1,2,3-cd)pyrene	27.14	276	5238	0.322	ng/ul#	74
25) Dibenzo(a,h)anthracene	27.14	278	1371	0.101	ng/ul#	79
26) Benzo(a,h,i)perylene	28.00	276	4407	0.311	ng/ul	96

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

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