

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE091218\
 Data File : BE097472.D
 Acq On : 13 Sep 2018 4:18
 Operator : SJ/JU
 Sample : J4792-19DL 5X
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A3YZ5DL

Manual Integrations
 APPROVED

Sohil
 9/13/2018 5:42:40 PM

Quant Time: Sep 13 08:08:02 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE082918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 13 07:48:38 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	1039	0.40	ng/ul	0.00
2) Naphthalene-d8	10.12	136	6127	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.01	164	4189	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.76	188	10202m	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	11746m	0.40	ng/ul	0.00
20) Perylene-d12	23.22	264	11457m	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	603	0.06	ng/ul	0.00
14) Fluoranthene-d10	18.80	212	2011m	0.06	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.17	128	2548	0.171	ng/ul#	95
5) 2-Methylnaphthalene	11.80	142	514	0.054	ng/ul	99
7) Acenaphthylene	13.73	152	553	0.036	ng/ul#	47
8) Acenaphthene	14.07	153	458	0.035	ng/ul#	90
9) Fluorene	15.07	166	874	0.058	ng/ul#	76
12) Phenanthrene	16.80	178	8981m	0.342	ng/ul	
13) Anthracene	16.89	178	2669m	0.121	ng/ul	
15) Fluoranthene	18.84	202	22594m	0.670	ng/ul	
17) Pyrene	19.20	202	30813m	0.975	ng/ul	
18) Benzo(a)anthracene	20.96	228	19384m	0.642	ng/ul	
19) Chrysene	21.01	228	23489m	0.676	ng/ul	
21) Benzo(b)fluoranthene	22.54	252	30154m	0.983	ng/ul	
22) Benzo(k)fluoranthene	22.58	252	8788m	0.257	ng/ul	
23) Benzo(a)pyrene	23.11	252	15561m	0.548	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.51	276	12463m	0.339	ng/ul	
25) Dibenzo(a,h)anthracene	25.52	278	4051m	0.130	ng/ul	
26) Benzo(g,h,i)perylene	26.22	276	11155m	0.356	ng/ul	

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

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