

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE122818\
 Data File : BE098591.D
 Acq On : 29 Dec 2018 1:26
 Operator : SJ/JU
 Sample : J6489-09DL 5X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F9F86DL

Manual Integrations
 APPROVED

Sohil
 1/2/2019 3:57:21 PM

Quant Time: Dec 29 03:56:31 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE122818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Dec 29 02:46:25 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.81	152	976	0.40	ng/ul	0.00
2) Naphthalene-d8	10.60	136	4451	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	3114	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	7304m	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	7982m	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	7558m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.20	152	420	0.06	ng/ul	0.00
14) Fluoranthene-d10	19.26	212	1264m	0.05	ng/ul	0.00
Target Compounds						
3) Naphthalene	10.65	128	375	0.031	ng/ul#	1
7) Acenaphthylene	14.19	152	324	0.024	ng/ul#	1
8) Acenaphthene	14.53	153	1505	0.143	ng/ul	96
9) Fluorene	15.52	166	13797	1.106	ng/ul	97
12) Phenanthrene	17.27	178	307362m	15.656	ng/ul	
13) Anthracene	17.36	178	13321m	0.810	ng/ul	
15) Fluoranthene	19.29	202	36476m	1.440	ng/ul	
17) Pyrene	19.65	202	95499	3.742	ng/ul	98
18) Benzo(a)anthracene	21.39	228	35956m	1.543	ng/ul	
19) Chrysene	21.45	228	62739m	2.173	ng/ul	
21) Benzo(b)fluoranthene	23.15	252	12957m	0.560	ng/ul	
22) Benzo(k)fluoranthene	23.20	252	2497m	0.098	ng/ul	
23) Benzo(a)pyrene	23.80	252	10479m	0.465	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.56	276	3944m	0.152	ng/ul	
25) Dibenzo(a,h)anthracene	26.58	278	2447m	0.116	ng/ul	
26) Benzo(a,h,i)perylene	27.37	276	5420m	0.241	ng/ul	

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

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