

Data Path : U:\HPCHEM1\BNA E\DATA\BE020218\
 Data File : BE095428.D
 Acq On : 3 Feb 2018 9:38
 Operator : SJ/JU
 Sample : J1295-13DL 5X
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6C11DL

Manual Integrations
 APPROVED

Sohil
 2/5/2018 5:49:07 PM

Quant Time: Feb 05 01:58:57 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 05 01:47:56 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.43	152	1982	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	6198	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	3704	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	7997	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	7894	0.40	ng/ul	0.00
20) Perylene-d12	24.43	264	8230	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	787	0.08	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	1948	0.06	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.74	152	645	0.036	ng/ul#	65
11) Pentachlorophenol	17.36	266	70540	45.541	ng/ul	97
12) Phenanthrene	17.75	178	7987	0.297	ng/ul#	91
13) Anthracene	17.84	178	1416	0.058	ng/ul	94
15) Fluoranthene	19.72	202	17532	0.503	ng/ul	84
17) Pyrene	20.07	202	15708	0.593	ng/ul#	80
18) Benzo(a)anthracene	21.78	228	2816	0.114	ng/ul#	89
19) Chrysene	21.84	228	7054	0.275	ng/ul	97
21) Benzo(b)fluoranthene	23.62	252	8038m	0.277	ng/ul	
22) Benzo(k)fluoranthene	23.66	252	2092m	0.079	ng/ul	
23) Benzo(a)pyrene	24.31	252	839	0.032	ng/ul#	58
24) Indeno(1,2,3-cd)pyrene	27.24	276	2120	0.075	ng/ul#	74
25) Dibenzo(a,h)anthracene	27.27	278	543	0.023	ng/ul#	40
26) Benzo(g,h,i)perylene	28.12	276	1675	0.070	ng/ul#	88

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

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