

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE020518\
 Data File : BE095496.D
 Acq On : 5 Feb 2018 18:25
 Operator : SJ/JU
 Sample : J1296-09
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6C27

Manual Integrations
 APPROVED

Sohil
 2/6/2018 2:30:55 PM

Quant Time: Feb 06 02:52:46 2018
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 06 02:09:13 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.43	152	1751	0.40	ng/ul	0.00
2) Naphthalene-d8	11.25	136	4805	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.01	164	3078	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	6833	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	7000	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	6061m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	2971	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	8529	0.33	ng/ul	0.00
Target Compounds						Qvalue
7) Acenaphthylene	14.74	152	1321	0.089	ng/ul#	90
11) Pentachlorophenol	17.36	266	10361	7.829	ng/ul	98
12) Phenanthrene	17.75	178	5432m	0.236	ng/ul	
13) Anthracene	17.85	178	2595m	0.123	ng/ul	
15) Fluoranthene	19.72	202	34932	1.174	ng/ul	85
17) Pyrene	20.07	202	32231m	1.372	ng/ul	
18) Benzo(a)anthracene	21.78	228	11892	0.543	ng/ul#	85
19) Chrysene	21.84	228	12284	0.541	ng/ul	98
21) Benzo(b)fluoranthene	23.62	252	17095m	0.799	ng/ul	
22) Benzo(k)fluoranthene	23.66	252	4447m	0.227	ng/ul	
23) Benzo(a)pyrene	24.32	252	4084m	0.214	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.23	276	4793m	0.229	ng/ul	
25) Dibenzo(a,h)anthracene	27.25	278	1076m	0.063	ng/ul	
26) Benzo(g,h,i)perylene	28.10	276	4292m	0.242	ng/ul	

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

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