

Data Path : Z:\HPCHEM1\BNA E\DATA\BE020118\
 Data File : BE095385.D
 Acq On : 2 Feb 2018 7:51
 Operator : SJ/JU
 Sample : J1295-14
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6C12

Manual Integrations
 APPROVED

Sohil
 2/2/2018 4:42:49 PM

Quant Time: Feb 02 08:33:02 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 02 04:53:51 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	2394	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	7483	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	4750	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	9587	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	10078	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	9075m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.82	152	3173	0.26	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	8815	0.24	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.74	152	25739	1.126	ng/ul	97
8) Acenaphthene	15.08	153	372	0.021	ng/ul#	65
9) Fluorene	16.02	166	2316	0.127	ng/ul#	70
11) Pentachlorophenol	17.36	266	33074	17.812	ng/ul	96
12) Phenanthrene	17.75	178	7891	0.245	ng/ul#	92
13) Anthracene	17.84	178	26939	0.913	ng/ul#	93
15) Fluoranthene	19.73	202	855992	20.503	ng/ul	81
17) Pyrene	20.08	202	1181945	34.951	ng/ul#	78
18) Benzo(a)anthracene	21.79	228	294985	9.362	ng/ul#	85
19) Chrysene	21.84	228	395809	12.097	ng/ul	95
21) Benzo(b)fluoranthene	23.62	252	400189m	12.499	ng/ul	
22) Benzo(k)fluoranthene	23.67	252	115035m	3.928	ng/ul	
23) Benzo(a)pyrene	24.32	252	150186m	5.254	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.25	276	90756m	2.895	ng/ul	
25) Dibenzo(a,h)anthracene	27.26	278	22531m	0.879	ng/ul	
26) Benzo(a,h,i)perylene	28.12	276	60122m	2.265	ng/ul	

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

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