

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE020918\
 Data File : BE095627.D
 Acq On : 10 Feb 2018 2:59
 Operator : SJ/JU
 Sample : J1339-07DL 5X
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6D40DL

Manual Integrations
 APPROVED

Sohil
 2/12/2018 11:39:34 AM

Quant Time: Feb 10 03:44:55 2018
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Feb 10 00:56:50 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.43	152	1734	0.40	ng/ul	0.00
2) Naphthalene-d8	11.25	136	4969	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.01	164	3116	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	7537	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	7336	0.40	ng/ul	0.00
20) Perylene-d12	24.43	264	7361	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	438	0.05	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	1245	0.04	ng/ul	0.00
Target Compounds						
					Qvalue	
9) Fluorene	16.03	166	492	0.041	ng/ul#	91
11) Pentachlorophenol	17.36	266	10598	7.260	ng/ul	98
12) Phenanthrene	17.75	178	12187m	0.481	ng/ul	
13) Anthracene	17.85	178	1414	0.061	ng/ul	96
15) Fluoranthene	19.72	202	12872	0.392	ng/ul	83
17) Pyrene	20.07	202	9635m	0.391	ng/ul	
18) Benzo(a)anthracene	21.78	228	2525	0.110	ng/ul#	87
19) Chrysene	21.84	228	2070m	0.087	ng/ul	
21) Benzo(b)fluoranthene	23.62	252	2059m	0.079	ng/ul	
22) Benzo(k)fluoranthene	23.66	252	517m	0.022	ng/ul	
23) Benzo(a)pyrene	24.31	252	816	0.035	ng/ul#	86
26) Benzo(g,h,i)perylene	28.11	276	431	0.020	ng/ul#	57

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

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