

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE020918\
 Data File : BE095629.D
 Acq On : 10 Feb 2018 4:11
 Operator : SJ/JU
 Sample : J1339-03DL 5X
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6C93DL

Manual Integrations
 APPROVED

Sohil
 2/12/2018 11:39:42 AM

Quant Time: Feb 12 00:38:13 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 02:59:57 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.43	152	1644	0.40	ng/ul	0.00
2) Naphthalene-d8	11.25	136	4650	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.01	164	2917	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	7233	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	6833	0.40	ng/ul	0.00
20) Perylene-d12	24.43	264	6668	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	212	0.03	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	588	0.02	ng/ul	0.00
Target Compounds				Qvalue		
11) Pentachlorophenol	17.36	266	1250	0.892	ng/ul	99
12) Phenanthrene	17.75	178	834	0.034	ng/ul#	88
15) Fluoranthene	19.72	202	4559	0.145	ng/ul	84
17) Pyrene	20.07	202	4555m	0.199	ng/ul	
18) Benzo(a)anthracene	21.78	228	2293	0.107	ng/ul#	88
19) Chrysene	21.83	228	1884m	0.085	ng/ul	
21) Benzo(b)fluoranthene	23.62	252	2727m	0.116	ng/ul	
22) Benzo(k)fluoranthene	23.66	252	664m	0.031	ng/ul	
23) Benzo(a)pyrene	24.31	252	977	0.047	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	27.23	276	679	0.029	ng/ul#	78
26) Benzo(g,h,i)perylene	28.10	276	576	0.030	ng/ul#	60

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

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