

Data Path : Z:\HPCHEM1\BNA E\DATA\BE012718\
 Data File : BE095197.D
 Acq On : 28 Jan 2018 9:18
 Operator : SJ/JU
 Sample : J1174-16DL 10X
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6A51DL

Manual Integrations
 APPROVED

Sohil
 1/29/2018 6:22:43 PM

Quant Time: Jan 29 01:29:56 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 Jan 29 01:25:04 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	2965	0.40	ng/ul	0.00
2) Naphthalene-d8	11.27	136	9481	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	5941	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	13193	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	14063	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	13341	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.82	152	380	0.02	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	1080	0.02	ng/ul	0.00
Target Compounds				Ovalue		
5) 2-Methylnaphthalene	12.89	142	5675	0.320	ng/ul	100
8) Acenaphthene	15.08	153	10448	0.471	ng/ul	97
9) Fluorene	16.03	166	9966	0.436	ng/ul#	74
11) Pentachlorophenol	17.36	266	113	0.044	ng/ul	97
12) Phenanthrene	17.75	178	63154	1.423	ng/ul#	87
13) Anthracene	17.85	178	9119	0.225	ng/ul#	92
15) Fluoranthene	19.72	202	51679	0.900	ng/ul	84
17) Pyrene	20.08	202	39172	0.830	ng/ul#	79
18) Benzo(a)anthracene	21.79	228	8212	0.187	ng/ul#	88
19) Chrysene	21.84	228	7193	0.158	ng/ul	95
21) Benzo(b)fluoranthene	23.62	252	5333m	0.113	ng/ul	
22) Benzo(k)fluoranthene	23.67	252	1922m	0.045	ng/ul	
23) Benzo(a)pyrene	24.32	252	3180	0.076	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	27.24	276	1349	0.029	ng/ul#	79
26) Benzo(a,h,i)perylene	28.11	276	1231	0.032	ng/ul#	77

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

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