

Data Path : U:\HPCHEM1\BNA E\DATA\BE012518\
 Data File : BE095115.D
 Acq On : 25 Jan 2018 14:15
 Operator : SJ/JU
 Sample : J1236-06DL 5X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 TW-2DL

Manual Integrations
 APPROVED

Sohil
 1/26/2018 11:21:59 AM

Quant Time: Jan 25 20:22:29 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE012318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 23 15:47:25 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.46	152	5381	0.40	ng	0.00
7) Naphthalene-d8	11.29	136	12121	0.40	ng	-0.02
13) Acenaphthene-d10	15.03	164	7482	0.40	ng	0.00
19) Phenanthrene-d10	17.74	188	14963	0.40	ng	0.00
27) Chrysene-d12	21.84	240	15246	0.40	ng	0.00
34) Perylene-d12	24.48	264	14184	0.40	ng	-0.01

System Monitoring Compounds						
4) 2-Fluorophenol	5.94	112	617	0.07	ng	0.00
5) Phenol-d6	7.59	99	3697	0.28	ng	0.01
8) Nitrobenzene-d5	9.63	82	5662	0.51	ng	0.00
11) 2-Methylnaphthalene-d10	12.83	152	1536	0.08	ng	0.00
14) 2,4,6-Tribromophenol	16.50	330	1199	0.57	ng	0.00
15) 2-Fluorobiphenyl	13.68	172	24847	0.76	ng	0.00
25) Fluoranthene-d10	19.72	212	16199	0.09	ng	0.00
29) Terphenyl-d14	20.28	244	145730	4.67	ng	0.00

Target Compounds				Qvalue		
9) Naphthalene	11.34	128	3117794	22.478	ng	99
12) 2-Methylnaphthalene	12.91	142	39057	1.643	ng	99
16) Acenaphthylene	14.76	152	917	0.023	ng	# 62
17) Acenaphthene	15.09	154	3545	0.136	ng	97
18) Fluorene	16.05	166	6910	0.212	ng	# 80
23) Phenanthrene	17.77	178	39840	0.862	ng	99
24) Anthracene	17.86	178	7681	0.186	ng	97
26) Fluoranthene	19.75	202	46049	0.829	ng	# 97
28) Pyrene	20.09	202	45376	0.704	ng	# 90
30) Benzo(a)anthracene	21.81	228	18779	0.394	ng	98
31) Chrysene	21.87	228	16148	0.335	ng	96
32) Bis(2-ethylhexyl)phthalate	21.70	149	257662	3.885	ng	100
33) Indeno(1,2,3-cd)pyrene	27.30	276	9136	0.208	ng	# 91
35) Benzo(b)fluoranthene	23.66	252	19314	0.384	ng	# 92
36) Benzo(k)fluoranthene	23.71	252	7096m	0.138	ng	
37) Benzo(a)pyrene	24.36	252	11571	0.251	ng	# 92
38) Dibenzo(a,h)anthracene	27.31	278	2367	0.061	ng	# 79
39) Benzo(a,h,i)perylene	28.18	276	9137	0.217	ng	# 91

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

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