

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA E\DATA\BE030718\
 Data File : BE095721.D
 Acq On : 7 Mar 2018 13:30
 Operator : SJ/JU
 Sample : J1811-10
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 T-10_030518

Manual Integrations
 APPROVED

Sohil
 3/8/2018 4:46:30 PM

Quant Time: Mar 07 14:35:25 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE022318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 23 12:12:42 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.43	152	1624	0.40	ng	0.00
7) Naphthalene-d8	11.26	136	6374	0.40	ng	-0.02
13) Acenaphthene-d10	15.00	164	3625	0.40	ng	0.00
19) Phenanthrene-d10	17.71	188	8106	0.40	ng	0.00
27) Chrysene-d12	21.80	240	8370	0.40	ng	0.00
34) Perylene-d12	24.43	264	8115	0.40	ng	0.00

System Monitoring Compounds						
4) 2-Fluorophenol	5.92	112	1043	0.24	ng	0.00
5) Phenol-d6	7.56	99	1465	0.25	ng	0.00
8) Nitrobenzene-d5	9.61	82	1585	0.24	ng	0.00
11) 2-Methylnaphthalene-d10	12.80	152	2390	0.22	ng	0.00
14) 2,4,6-Tribromophenol	16.47	330	338	0.31	ng	0.00
15) 2-Fluorobiphenyl	13.64	172	3364	0.21	ng	0.00
25) Fluoranthene-d10	19.69	212	20458	0.19	ng	0.00
29) Terphenyl-d14	20.25	244	3194	0.19	ng	0.00

Target Compounds				Qvalue		
9) Naphthalene	11.31	128	1874	0.026	ng	# 91
12) 2-Methylnaphthalene	12.88	142	677	0.053	ng	95
16) Acenaphthylene	14.74	152	1881	0.092	ng	96
18) Fluorene	16.03	166	1722	0.109	ng	# 77
23) Phenanthrene	17.75	178	40715	1.607	ng	98
24) Anthracene	17.84	178	8341	0.359	ng	95
26) Fluoranthene	19.72	202	72476	2.325	ng	# 97
28) Pyrene	20.07	202	70114	2.048	ng	96
30) Benzo(a)anthracene	21.78	228	28284	1.036	ng	99
31) Chrysene	21.84	228	26169	0.964	ng	95
32) Bis(2-ethylhexyl)phthalate	21.66	149	4111	0.085	ng	# 95
33) Indeno(1,2,3-cd)pyrene	27.23	276	9822	0.373	ng	# 91
35) Benzo(b)fluoranthene	23.61	252	24351	0.842	ng	# 91
36) Benzo(k)fluoranthene	23.66	252	9253m	0.322	ng	
37) Benzo(a)pyrene	24.30	252	16816	0.645	ng	# 92
38) Dibenzo(a,h)anthracene	27.24	278	2864	0.124	ng	# 84
39) Benzo(a,h,i)perylene	28.11	276	9098	0.366	ng	# 91

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

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