

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE052319\
 Data File : BE099732.D
 Acq On : 23 May 2019 13:55
 Operator : JU/SJ
 Sample : K2966-06DL 20X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 FESB-18(2.5-3.0)DL

Manual Integrations
 APPROVED

Sohil
 5/24/2019 5:02:41 PM

Quant Time: May 23 15:15:46 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE052019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon May 20 19:29:40 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.80	152	1167	0.40	ng	-0.01
7) Naphthalene-d8	10.61	136	4180	0.40	ng	0.00
13) Acenaphthene-d10	14.47	164	2918	0.40	ng	-0.01
19) Phenanthrene-d10	17.22	188	8443	0.40	ng	-0.01
27) Chrysene-d12	21.40	240	12780	0.40	ng	-0.01
34) Perylene-d12	23.94	264	17000	0.40	ng	-0.02
System Monitoring Compounds						
4) 2-Fluorophenol	5.36	112	65	0.02	ng	-0.01
5) Phenol-d6	6.96	99	128	0.03	ng	-0.01
8) Nitrobenzene-d5	8.96	82	86	0.02	ng	-0.01
11) 2-Methylnaphthalene-d10	12.21	152	89	0.01	ng	-0.01
14) 2,4,6-Tribromophenol	15.97	330	44	0.16	ng	-0.01
15) 2-Fluorobiphenyl	13.10	172	136	0.01	ng	0.00
25) Fluoranthene-d10	19.25	212	805	0.01	ng	-0.01
29) Terphenyl-d14	19.85	244	260	0.01	ng	-0.02
Target Compounds						Qvalue
9) Naphthalene	10.65	128	2669	0.060	ng	# 91
12) 2-Methylnaphthalene	12.28	142	308	0.042	ng	# 86
16) Acenaphthylene	14.20	152	2600	0.191	ng	98
17) Acenaphthene	14.54	154	1204	0.158	ng	99
18) Fluorene	15.52	166	2371	0.211	ng	94
22) Pentachlorophenol	16.89	266	80	0.234	ng	91
23) Phenanthrene	17.26	178	46629	2.260	ng	100
24) Anthracene	17.35	178	7366	0.394	ng	98
26) Fluoranthene	19.29	202	123271	3.779	ng	99
28) Pvrene	19.65	202	117688	3.665	ng	# 95
30) Benzo(a)anthracene	21.39	228	82415	2.269	ng	98
31) Chrysene	21.44	228	71115	1.832	ng	98
32) Bis(2-ethylhexyl)phthalate	21.31	149	8699	0.208	ng	99
33) Indeno(1,2,3-cd)pyrene	26.62	276	50717	1.082	ng	98
35) Benzo(b)fluoranthene	23.16	252	94795	2.035	ng	99
36) Benzo(k)fluoranthene	23.21	252	38921m	0.802	ng	
37) Benzo(a)pvrene	23.82	252	73499	1.640	ng	99
38) Dibenzo(a,h)anthracene	26.64	278	13255	0.282	ng	# 90
39) Benzo(g,h,i)perylene	27.46	276	50218	1.056	ng	99

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

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