

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE110719\
 Data File : BE100708.D
 Acq On : 7 Nov 2019 19:13
 Operator : JU
 Sample : K5725-04
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SO-4

Manual Integrations
 APPROVED

Sohil
 11/8/2019 3:32:30 PM

Quant Time: Nov 08 11:40:13 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE101019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 23 16:49:33 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.85	152	2559	0.40	ng	0.00
7) Naphthalene-d8	10.64	136	13884	0.40	ng	0.01
13) Acenaphthene-d10	14.50	164	25100	0.40	ng	0.03
19) Phenanthrene-d10	17.23	188	15772	0.40	ng	0.03
27) Chrysene-d12	21.39	240	21477	0.40	ng	0.04
34) Perylene-d12	23.88	264	28380	0.40	ng	0.05

System Monitoring Compounds						
4) 2-Fluorophenol	5.46	112	2134	0.27	ng	0.01
5) Phenol-d6	7.02	99	3729	0.33	ng	0.01
8) Nitrobenzene-d5	9.00	82	9352m	1.09	ng	0.01
11) 2-Methylnaphthalene-d10	12.24	152	12580	0.57	ng	0.01
14) 2,4,6-Tribromophenol	15.98	330	1004	0.15	ng	0.04
15) 2-Fluorobiphenyl	13.12	172	7818	0.09	ng	0.01
25) Fluoranthene-d10	19.24	212	113999	0.56	ng	0.02
29) Terphenyl-d14	19.83	244	18525	0.39	ng	0.02

Target Compounds						Qvalue
9) Naphthalene	10.69	128	1823379	12.489	ng	# 99
12) 2-Methylnaphthalene	12.31	142	531959	21.376	ng	# 90
16) Acenaphthylene	14.21	152	375724	3.479	ng	# 69
17) Acenaphthene	14.56	154	188675	2.703	ng	# 68
18) Fluorene	15.53	166	335519	3.619	ng	# 64
23) Phenanthrene	17.27	178	2537056	58.482	ng	# 93
24) Anthracene	17.36	178	646465	16.060	ng	# 81
26) Fluoranthene	19.31	202	4854194	85.732	ng	# 96
28) Pyrene	19.65	202	4922998	82.045	ng	# 88
30) Benzo(a)anthracene	21.37	228	3192342	45.567	ng	# 92
31) Chrysene	21.43	228	3089056m	44.120	ng	# 92
32) Bis(2-ethylhexyl)phthalate	21.29	149	453597	3.095	ng	# 98
33) Indeno(1,2,3-cd)pyrene	26.55	276	3414394	39.679	ng	# 96
35) Benzo(b)fluoranthene	23.13	252	4934225	59.626	ng	# 98
36) Benzo(k)fluoranthene	23.17	252	1895044m	22.266	ng	# 98
37) Benzo(a)pyrene	23.78	252	3604428	46.793	ng	# 98
38) Dibenzo(a,h)anthracene	26.55	278	789090	9.911	ng	# 95
39) Benzo(a,h,i)perylene	27.37	276	3367490	42.054	ng	# 98

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

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