

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE042619\
 Data File : BE099394.D
 Acq On : 27 Apr 2019 1:21
 Operator : JU/SJ
 Sample : K2533-19
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 EXT-027-SW019-042219

Manual Integrations
 APPROVED

Sohil
 4/29/2019 1:31:58 PM

Quant Time: Apr 27 07:14:11 2019

Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE041819.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Fri Apr 19 06:51:38 2019

Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.82	152	3400	0.40	ng	0.00
7) Naphthalene-d8	10.61	136	13629	0.40	ng	0.01
13) Acenaphthene-d10	14.46	164	10282	0.40	ng	0.00
19) Phenanthrene-d10	17.20	188	28113	0.40	ng	0.01
27) Chrysene-d12	21.37	240	31955	0.40	ng	0.00
34) Perylene-d12	23.88	264	36677	0.40	ng	0.02

System Monitoring Compounds

4) 2-Fluorophenol	5.40	112	3662	0.44	ng	0.01
5) Phenol-d6	6.97	99	5267	0.44	ng	0.00
8) Nitrobenzene-d5	8.98	82	4797	0.45	ng	0.01
11) 2-Methylnaphthalene-d10	12.21	152	9338	0.39	ng	0.01
14) 2,4,6-Tribromophenol	15.94	330	2521	0.74	ng	0.00
15) 2-Fluorobiphenyl	13.09	172	13731	0.34	ng	0.00
25) Fluoranthene-d10	19.22	212	126392	0.37	ng	0.00
29) Terphenyl-d14	19.82	244	21563	0.34	ng	0.00

Target Compounds

						Qvalue
9) Naphthalene	10.65	128	5476	0.037	ng	# 95
12) 2-Methylnaphthalene	12.28	142	833	0.033	ng	# 95
16) Acenaphthylene	14.18	152	4345	0.093	ng	98
17) Acenaphthene	14.53	154	582	0.020	ng	93
18) Fluorene	15.50	166	2343	0.057	ng	# 88
23) Phenanthrene	17.24	178	47292	0.635	ng	99
24) Anthracene	17.32	178	4050	0.061	ng	# 78
26) Fluoranthene	19.25	202	82449	0.827	ng	98
28) Pyrene	19.62	202	85374	0.907	ng	# 96
30) Benzo(a)anthracene	21.35	228	48014	0.511	ng	# 94
31) Chrysene	21.40	228	50592m	0.506	ng	
32) Bis(2-ethylhexyl)phthalate	21.27	149	48842	0.571	ng	# 99
33) Indeno(1,2,3-cd)pyrene	26.52	276	20907	0.212	ng	95
35) Benzo(b)fluoranthene	23.10	252	54930	0.489	ng	# 76
36) Benzo(k)fluoranthene	23.15	252	21663m	0.194	ng	
37) Benzo(a)pyrene	23.76	252	39444	0.392	ng	# 86
38) Dibenzo(a,h)anthracene	26.54	278	5544	0.054	ng	# 25
39) Benzo(a,h,i)perylene	27.35	276	18935	0.182	ng	# 82

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

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