

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE120519\
 Data File : BE100793.D
 Acq On : 5 Dec 2019 17:39
 Operator : JU
 Sample : K6053-05
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 EXT-029-SW180-112519

Manual Integrations
 APPROVED

mohammad
 12/6/2019 3:02:43 PM

Quant Time: Dec 05 18:20:37 2019

Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE120419.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Wed Dec 04 17:18:11 2019

Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.76	152	6044	0.40	ng	-0.01
7) Naphthalene-d8	10.54	136	27905	0.40	ng	0.00
13) Acenaphthene-d10	14.40	164	16705	0.40	ng	0.00
19) Phenanthrene-d10	17.12	188	40805	0.40	ng	-0.01
27) Chrysene-d12	21.30	240	42349	0.40	ng	0.00
34) Perylene-d12	23.76	264	49908	0.40	ng	0.00

System Monitoring Compounds

4) 2-Fluorophenol	5.38	112	5417	0.38	ng	0.00
5) Phenol-d6	6.93	99	9200	0.42	ng	0.00
8) Nitrobenzene-d5	8.90	82	5997	0.45	ng	0.00
11) 2-Methylnaphthalene-d10	12.14	152	10748	0.27	ng	0.00
14) 2,4,6-Tribromophenol	15.88	330	1468	0.79	ng	0.00
15) 2-Fluorobiphenyl	13.03	172	15044	0.23	ng	0.00
25) Fluoranthene-d10	19.16	212	121489	0.25	ng	0.00
29) Terphenyl-d14	19.77	244	23463	0.23	ng	0.00

Target Compounds

						Qvalue
9) Naphthalene	10.59	128	27185	0.092	ng	99
12) 2-Methylnaphthalene	12.21	142	6359	0.137	ng	97
16) Acenaphthylene	14.11	152	31033	0.449	ng	99
17) Acenaphthene	14.46	154	4049	0.085	ng	97
18) Fluorene	15.43	166	30110m	0.490	ng	
23) Phenanthrene	17.16	178	423763	3.429	ng	100
24) Anthracene	17.25	178	101183	0.979	ng	# 92
26) Fluoranthene	19.19	202	484338	3.283	ng	99
28) Pyrene	19.55	202	412461	2.998	ng	# 95
30) Benzo(a)anthracene	21.29	228	284362	2.278	ng	94
31) Chrysene	21.34	228	232464m	1.585	ng	
32) Bis(2-ethylhexyl)phthalate	21.24	149	69991	0.769	ng	# 94
33) Indeno(1,2,3-cd)pyrene	26.33	276	78051	0.611	ng	98
35) Benzo(b)fluoranthene	23.01	252	229567	1.680	ng	96
36) Benzo(k)fluoranthene	23.05	252	83891m	0.547	ng	
37) Benzo(a)pyrene	23.64	252	166875	1.411	ng	98
38) Dibenzo(a,h)anthracene	26.35	278	26854	0.228	ng	# 78
39) Benzo(a,h,i)perylene	27.12	276	67871	0.539	ng	96

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

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