

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE101519\
 Data File : BE100648.D
 Acq On : 15 Oct 2019 14:11
 Operator : JU
 Sample : K4994-02DL 10X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 EXT-SW091-091819-3DL

Manual Integrations
 APPROVED

mohammad
 10/17/2019 7:56:52 AM

Quant Time: Oct 15 16:32:06 2019

Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE101019.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Thu Oct 10 14:01:52 2019

Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.86	152	4919	0.40	ng	0.00
7) Naphthalene-d8	10.64	136	22454	0.40	ng	0.00
13) Acenaphthene-d10	14.49	164	14425	0.40	ng	0.00
19) Phenanthrene-d10	17.20	188	29288	0.40	ng	0.00
27) Chrysene-d12	21.38	240	31209	0.40	ng	0.01
34) Perylene-d12	23.86	264	35676	0.40	ng	0.00

System Monitoring Compounds

4) 2-Fluorophenol	5.46	112	352	0.02	ng	0.01
5) Phenol-d6	7.02	99	576	0.03	ng	0.01
8) Nitrobenzene-d5	9.00	82	489	0.04	ng	0.01
11) 2-Methylnaphthalene-d10	12.24	152	834	0.02	ng	0.01
14) 2,4,6-Tribromophenol	15.96	330	129	0.03	ng	0.00
15) 2-Fluorobiphenyl	13.12	172	1029	0.02	ng	0.00
25) Fluoranthene-d10	19.23	212	6986	0.02	ng	0.00
29) Terphenyl-d14	19.83	244	2059	0.03	ng	0.00

Target Compounds

						Qvalue
9) Naphthalene	10.69	128	78064	0.331	ng	100
12) 2-Methylnaphthalene	12.31	142	9125	0.227	ng	96
16) Acenaphthylene	14.20	152	47662	0.768	ng	98
17) Acenaphthene	14.54	154	6430	0.160	ng	97
18) Fluorene	15.51	166	27150	0.510	ng	87
23) Phenanthrene	17.25	178	485187	6.023	ng	99
24) Anthracene	17.34	178	46390	0.621	ng	99
26) Fluoranthene	19.27	202	714955	6.800	ng	99
28) Pyrene	19.62	202	830786	9.528	ng	# 96
30) Benzo(a)anthracene	21.35	228	400871m	3.938	ng	
31) Chrysene	21.41	228	485105	4.768	ng	97
32) Bis(2-ethylhexyl)phthalate	21.30	149	8909	0.042	ng	# 80
33) Indeno(1,2,3-cd)pyrene	26.48	276	231578	1.852	ng	96
35) Benzo(b)fluoranthene	23.10	252	476690	4.582	ng	98
36) Benzo(k)fluoranthene	23.15	252	166930m	1.560	ng	
37) Benzo(a)pyrene	23.75	252	370284	3.824	ng	98
38) Dibenzo(a,h)anthracene	26.51	278	65623	0.656	ng	# 83
39) Benzo(a,h,i)perylene	27.29	276	246666	2.450	ng	98

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

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