

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE110819\
 Data File : BE100717.D
 Acq On : 8 Nov 2019 14:57
 Operator : JU
 Sample : K5725-04DL2 50X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleID :
 SO-4DL2

Manual Integrations
 APPROVED

mohammad
 11/11/2019 8:09:55 AM

Quant Time: Nov 08 15:42:18 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	2231	0.40	ng	0.00
7) Naphthalene-d8	10.64	136	10841	0.40	ng	0.01
13) Acenaphthene-d10	14.47	164	7676	0.40	ng	0.00
19) Phenanthrene-d10	17.20	188	16166	0.40	ng	0.00
27) Chrysene-d12	21.37	240	17104	0.40	ng	0.01
34) Perylene-d12	23.85	264	19350	0.40	ng	0.01

System Monitoring Compounds						
4) 2-Fluorophenol	5.45	112	42	0.01	ng	0.01
5) Phenol-d6	7.03	99	70	0.01	ng	0.02
8) Nitrobenzene-d5	8.99	82	498	0.07	ng	0.00
11) 2-Methylnaphthalene-d10	12.24	152	222	0.01	ng	0.01
14) 2,4,6-Tribromophenol	15.96	330	9	0.00	ng	0.01
15) 2-Fluorobiphenyl	13.10	172	76	0.00	ng	0.00
25) Fluoranthene-d10	19.22	212	1221	0.01	ng	0.00
29) Terphenyl-d14	19.82	244	381	0.01	ng	0.00

Target Compounds						Qvalue
9) Naphthalene	10.68	128	33361	0.293	ng	# 98
12) 2-Methylnaphthalene	12.30	142	10187	0.524	ng	# 89
16) Acenaphthylene	14.20	152	7928	0.240	ng	# 79
17) Acenaphthene	14.54	154	3966	0.186	ng	# 64
18) Fluorene	15.51	166	6846	0.241	ng	# 62
23) Phenanthrene	17.24	178	57219	1.287	ng	# 94
24) Anthracene	17.34	178	13420	0.325	ng	# 80
26) Fluoranthene	19.26	202	115776	1.995	ng	100
28) Pyrene	19.62	202	120187	2.515	ng	# 94
30) Benzo(a)anthracene	21.34	228	67205	1.205	ng	96
31) Chrysene	21.40	228	64326	1.154	ng	97
32) Bis(2-ethylhexyl)phthalate	21.28	149	10962	0.094	ng	100
33) Indeno(1,2,3-cd)pyrene	26.47	276	61906	0.903	ng	96
35) Benzo(b)fluoranthene	23.09	252	107148	1.899	ng	99
36) Benzo(k)fluoranthene	23.13	252	37108m	0.639	ng	
37) Benzo(a)pyrene	23.74	252	74386	1.416	ng	99
38) Dibenzo(a,h)anthracene	26.48	278	14296	0.263	ng	# 86
39) Benzo(a,h,i)perylene	27.28	276	64766	1.186	ng	99

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

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