

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE101519\
 Data File : BE100646.D
 Acq On : 15 Oct 2019 12:14
 Operator : JU
 Sample : K4994-10
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PST-SW117-091919-3

Manual Integrations
 APPROVED

mohammad
 10/17/2019 7:56:49 AM

Quant Time: Oct 15 13:03:46 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	3392	0.40	ng	0.00
7) Naphthalene-d8	10.64	136	15157	0.40	ng	0.00
13) Acenaphthene-d10	14.48	164	10469	0.40	ng	0.00
19) Phenanthrene-d10	17.20	188	25624	0.40	ng	0.00
27) Chrysene-d12	21.37	240	30795	0.40	ng	0.00
34) Perylene-d12	23.85	264	35247	0.40	ng	0.00

System Monitoring Compounds						
4) 2-Fluorophenol	5.44	112	3397	0.32	ng	0.00
5) Phenol-d6	7.01	99	5015	0.34	ng	0.00
8) Nitrobenzene-d5	8.99	82	4074	0.43	ng	0.00
11) 2-Methylnaphthalene-d10	12.24	152	7669	0.32	ng	0.01
14) 2,4,6-Tribromophenol	15.96	330	1539	0.56	ng	0.00
15) 2-Fluorobiphenyl	13.12	172	9983	0.26	ng	0.00
25) Fluoranthene-d10	19.23	212	96057	0.29	ng	0.00
29) Terphenyl-d14	19.83	244	20228	0.30	ng	0.00

Target Compounds				Qvalue		
9) Naphthalene	10.69	128	60084	0.377	ng	100
12) 2-Methylnaphthalene	12.31	142	13194	0.486	ng	100
16) Acenaphthylene	14.20	152	47008	1.044	ng	98
17) Acenaphthene	14.54	154	30897	1.061	ng	98
18) Fluorene	15.51	166	50538	1.307	ng	88
23) Phenanthrene	17.24	178	825764	11.716	ng	100
24) Anthracene	17.34	178	142181	2.174	ng	# 90
26) Fluoranthene	19.26	202	1051909	11.435	ng	99
28) Pyrene	19.62	202	1063925	12.366	ng	# 94
30) Benzo(a)anthracene	21.36	228	594336	5.916	ng	93
31) Chrysene	21.42	228	591901m	5.896	ng	
32) Bis(2-ethylhexyl)phthalate	21.31	149	1243828	5.920	ng	# 99
33) Indeno(1,2,3-cd)pyrene	26.48	276	284472	2.306	ng	96
35) Benzo(b)fluoranthene	23.10	252	634965	6.178	ng	98
36) Benzo(k)fluoranthene	23.15	252	247214m	2.339	ng	
37) Benzo(a)pyrene	23.75	252	467751	4.889	ng	99
38) Dibenzo(a,h)anthracene	26.50	278	83214	0.842	ng	# 86
39) Benzo(a,h,i)perylene	27.29	276	282143	2.837	ng	99

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

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