

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE070119\
 Data File : BE100289.D
 Acq On : 2 Jul 2019 6:33
 Operator : JU/SJ
 Sample : K3445-06DL 5X
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PST-007-SW050-061819DL

Manual Integrations
 APPROVED

mohammad
 7/3/2019 10:53:11 PM

Quant Time: Jul 02 08:18:50 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE062819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 28 14:06:53 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	824	0.40	ng	0.00
7) Naphthalene-d8	10.45	136	2974	0.40	ng	0.00
13) Acenaphthene-d10	14.33	164	2247	0.40	ng	0.02
19) Phenanthrene-d10	17.07	188	6395	0.40	ng	0.01
27) Chrysene-d12	21.24	240	7968	0.40	ng	0.00
34) Perylene-d12	23.68	264	9194	0.40	ng	0.02

System Monitoring Compounds						
4) 2-Fluorophenol	5.22	112	141	0.07	ng	0.00
5) Phenol-d6	6.85	99	155	0.06	ng	0.00
8) Nitrobenzene-d5	8.83	82	172	0.06	ng	0.01
11) 2-Methylnaphthalene-d10	12.05	152	285	0.05	ng	0.00
14) 2,4,6-Tribromophenol	15.82	330	96	0.06	ng	0.00
15) 2-Fluorobiphenyl	12.94	172	552	0.06	ng	0.00
25) Fluoranthene-d10	19.10	212	4312	0.05	ng	0.00
29) Terphenyl-d14	19.70	244	1585	0.10	ng	0.00

Target Compounds						Qvalue
9) Naphthalene	10.48	128	1880	0.057	ng	# 90
12) 2-Methylnaphthalene	12.12	142	162	0.028	ng	# 81
16) Acenaphthylene	14.04	152	1744	0.162	ng	98
17) Acenaphthene	14.38	154	460	0.075	ng	93
18) Fluorene	15.37	166	986	0.109	ng	# 97
23) Phenanthrene	17.11	178	33331	2.150	ng	100
24) Anthracene	17.20	178	4620	0.331	ng	# 92
26) Fluoranthene	19.13	202	76415	3.095	ng	100
28) Pyrene	19.49	202	69983	3.317	ng	# 96
30) Benzo(a)anthracene	21.23	228	43076	1.644	ng	97
31) Chrysene	21.28	228	38107	1.449	ng	96
33) Indeno(1,2,3-cd)pyrene	26.24	276	21811	0.627	ng	# 95
35) Benzo(b)fluoranthene	22.94	252	48255	1.936	ng	# 98
36) Benzo(k)fluoranthene	22.98	252	15871m	0.570	ng	
37) Benzo(a)pyrene	23.57	252	35230m	1.410	ng	
38) Dibenzo(a,h)anthracene	26.26	278	5086	0.195	ng	# 69
39) Benzo(a,h,i)perylene	27.04	276	20652	0.770	ng	96

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

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