

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE092519\
 Data File : BE100558.D
 Acq On : 25 Sep 2019 16:02
 Operator : JU
 Sample : K4995-17
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 EXT-SW164-091819-1

Manual Integrations
 APPROVED

mohammad
 9/26/2019 3:55:43 PM

Quant Time: Sep 26 01:11:03 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE092019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 20 18:26:17 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.86	152	7597	0.40	ng	0.00
7) Naphthalene-d8	10.64	136	36431	0.40	ng	-0.01
13) Acenaphthene-d10	14.48	164	22376	0.40	ng	0.00
19) Phenanthrene-d10	17.20	188	51103	0.40	ng	0.00
27) Chrysene-d12	21.37	240	58672	0.40	ng	0.00
34) Perylene-d12	23.84	264	67238	0.40	ng	0.00

System Monitoring Compounds						
4) 2-Fluorophenol	5.43	112	6484	0.36	ng	0.00
5) Phenol-d6	7.01	99	9354	0.38	ng	-0.01
8) Nitrobenzene-d5	9.00	82	7336	0.35	ng	0.00
11) 2-Methylnaphthalene-d10	12.24	152	15839	0.29	ng	0.00
14) 2,4,6-Tribromophenol	15.96	330	2034	0.56	ng	0.00
15) 2-Fluorobiphenyl	13.12	172	19801	0.25	ng	0.00
25) Fluoranthene-d10	19.22	212	182998	0.30	ng	0.00
29) Terphenyl-d14	19.82	244	43531	0.31	ng	0.00

Target Compounds					Qvalue	
9) Naphthalene	10.69	128	82003	0.217	ng	100
12) 2-Methylnaphthalene	12.31	142	16445	0.274	ng	99
16) Acenaphthylene	14.21	152	69427	0.754	ng	99
17) Acenaphthene	14.54	154	4583	0.074	ng	93
18) Fluorene	15.51	166	13232	0.170	ng	# 84
22) Pentachlorophenol	16.85	266	864	0.349	ng	96
23) Phenanthrene	17.24	178	305191	2.143	ng	100
24) Anthracene	17.34	178	43722	0.365	ng	97
26) Fluoranthene	19.25	202	556708	3.242	ng	98
28) Pvrene	19.62	202	669061	3.963	ng	96
30) Benzo(a)anthracene	21.34	228	330137	2.107	ng	94
31) Chrysene	21.40	228	399470	2.237	ng	96
32) Bis(2-ethylhexyl)phthalate	21.28	149	604417	3.579	ng	100
33) Indeno(1,2,3-cd)pyrene	26.45	276	231960	1.169	ng	97
35) Benzo(b)fluoranthene	23.08	252	463738	2.549	ng	# 97
36) Benzo(k)fluoranthene	23.12	252	167086m	0.794	ng	
37) Benzo(a)pvrene	23.73	252	330167	1.945	ng	97
38) Dibenzo(a,h)anthracene	26.46	278	64665	0.366	ng	# 83
39) Benzo(g,h,i)perylene	27.25	276	249541	1.352	ng	96

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

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