

Data Path : Z:\HPCHEM1\BNA E\DATA\BE020117\
 Data File : BE092712.D
 Acq On : 1 Feb 2017 13:14
 Operator : SJ/MA
 Sample : PB96613BS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PB96613BS

Manual Integrations
 APPROVED

mohammad
 2/2/2017 10:14:24 AM

Quant Time: Feb 01 14:53:56 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE011617.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 17 10:14:01 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.12	152	12543	5.00	ng	-0.02
7) Naphthalene-d8	10.91	136	54192	5.00	ng	0.00
12) Acenaphthene-d10	14.78	164	32621	5.00	ng	0.00
18) Phenanthrene-d10	17.53	188	71930	5.00	ng	0.00
24) Chrysene-d12	21.72	240	71674	5.00	ng	0.00
31) Perylene-d12	24.06	264	78982	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Fluorophenol	5.64	112	27746	9.91	ng	-0.02
5) Phenol-d6	7.29	99	37476	10.92	ng	-0.04
8) Nitrobenzene-d5	9.29	82	16508	4.69	ng	-0.04
13) 2,4,6-Tribromophenol	16.27	330	9098	9.42	ng	-0.02
14) 2-Fluorobiphenyl	13.40	172	34908	4.35	ng	-0.02
26) Terphenyl-d14	20.14	244	46757	4.61	ng	-0.02

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.37	88	5669	3.34	ng	# 77
3) n-Nitrosodimethylamine	3.72	42	9758	4.31	ng	# 98
6) bis(2-Chloroethyl)ether	7.54	93	16906	4.64	ng	# 87
9) Naphthalene	10.95	128	48718	4.31	ng	# 94
10) Hexachlorobutadiene	11.24	225	7369	4.41	ng	# 99
11) 2-Methylnaphthalene	12.58	142	31634	4.50	ng	# 97
15) Acenaphthylene	14.49	152	205269	4.57	ng	# 100
16) Acenaphthene	14.83	154	30863	4.28	ng	# 93
17) Fluorene	15.82	166	40636	4.49	ng	# 99
19) Hexachlorobenzene	16.86	284	10335	4.39	ng	# 39
20) Pentachlorophenol	17.21	266	9540	12.90	ng	# 88
21) Phenanthrene	17.58	178	65808	4.14	ng	# 95
22) Anthracene	17.67	178	65101	4.44	ng	# 99
23) Fluoranthene	19.58	202	314204	4.46	ng	# 98
25) Pyrene	19.93	202	335352	4.75	ng	# 99
27) Benzo(a)anthracene	21.70	228	315903	4.47	ng	# 71
28) Chrysene	21.75	228	351330m	4.42	ng	# 70
29) Bis(2-ethylhexyl)phthalate	21.61	149	70433	5.06	ng	# 98
30) Indeno(1,2,3-cd)pyrene	26.53	276	404656	4.72	ng	# 97
32) Benzo(b)fluoranthene	23.34	252	93007	4.87	ng	# 94
33) Benzo(k)fluoranthene	23.38	252	85344	4.86	ng	# 96
34) Benzo(a)pyrene	23.95	252	87131	4.72	ng	# 97
35) Dibenzo(a,h)anthracene	26.55	278	82959	4.62	ng	# 97
36) Benzo(g,h,i)perylene	27.28	276	342686	4.62	ng	# 97

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

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