

Data Path : Z:\HPCHEM1\BNA E\DATA\BE011917\
 Data File : BE092662.D
 Acq On : 19 Jan 2017 18:10
 Operator : UM/SJ
 Sample : PB96141BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleID :
 PB96141BS

Manual Integrations
 APPROVED

mohammad
 1/20/2017 2:05:25 PM

Quant Time: Jan 20 02:12:18 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	8.14	152	9640	5.00	ng	0.00
7) Naphthalene-d8	10.91	136	43422	5.00	ng	0.00
12) Acenaphthene-d10	14.78	164	26465	5.00	ng	0.00
18) Phenanthrene-d10	17.53	188	56978	5.00	ng	0.00
24) Chrysene-d12	21.72	240	61897	5.00	ng	0.00
31) Perylene-d12	24.06	264	72482	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Fluorophenol	5.64	112	21299	9.90	ng	-0.01
5) Phenol-d6	7.31	99	26862	10.18	ng	-0.02
8) Nitrobenzene-d5	9.29	82	13524	4.80	ng	-0.05
13) 2,4,6-Tribromophenol	16.27	330	7376	9.42	ng	-0.02
14) 2-Fluorobiphenyl	13.40	172	28062	4.31	ng	-0.02
26) Terphenyl-d14	20.14	244	39662	4.53	ng	-0.02

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.38	88	4733	3.64	ng	# 70
3) n-Nitrosodimethylamine	3.72	42	7848	4.51	ng	# 93
6) bis(2-Chloroethyl)ether	7.54	93	12797	4.57	ng	# 94
9) Naphthalene	10.95	128	38192	4.21	ng	# 95
10) Hexachlorobutadiene	11.24	225	5538	4.14	ng	# 99
11) 2-Methylnaphthalene	12.59	142	25200m	4.48	ng	# 99
15) Acenaphthylene	14.49	152	165924	4.55	ng	# 100
16) Acenaphthene	14.85	154	24436	4.18	ng	# 93
17) Fluorene	15.82	166	32589	4.44	ng	# 99
19) Hexachlorobenzene	16.86	284	8593m	4.61	ng	# 99
20) Pentachlorophenol	17.21	266	9182	15.68	ng	# 86
21) Phenanthrene	17.58	178	57107	4.54	ng	# 94
22) Anthracene	17.67	178	56062	4.83	ng	# 98
23) Fluoranthene	19.58	202	258626	4.64	ng	# 98
25) Pyrene	19.93	202	279665	4.58	ng	# 99
27) Benzo(a)anthracene	21.68	228	276951	4.54	ng	# 71
28) Chrysene	21.75	228	293800m	4.28	ng	# 99
29) Bis(2-ethylhexyl)phthalate	21.61	149	50138	4.21	ng	# 69
30) Indeno(1,2,3-cd)pyrene	26.53	276	375250	5.06	ng	# 98
32) Benzo(b)fluoranthene	23.34	252	83578	4.77	ng	# 97
33) Benzo(k)fluoranthene	23.38	252	79531	4.52	ng	# 94
34) Benzo(a)pyrene	23.95	252	80564	4.89	ng	# 95
35) Dibenzo(a,h)anthracene	26.55	278	77064	4.78	ng	# 95
36) Benzo(g,h,i)perylene	27.28	276	320478	4.71	ng	# 98

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

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