

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE080517\
 Data File : BE093685.D
 Acq On : 5 Aug 2017 21:21
 Operator : SJ/JU
 Sample : I4427-08
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 LASB-14A-2

Quant Time: Aug 06 03:39:08 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE080117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 02 10:58:31 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	3524	0.40	ng	0.00
7) Naphthalene-d8	10.71	136	17417	0.40	ng	0.00
13) Acenaphthene-d10	14.53	164	9889	0.40	ng	0.00
19) Phenanthrene-d10	17.24	188	24799	0.40	ng	0.00
27) Chrysene-d12	21.40	240	24864	0.40	ng	-0.01
34) Perylene-d12	23.91	264	22283	0.40	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Fluorophenol	5.52	112	3179	0.30	ng	0.00
5) Phenol-d6	7.09	99	4954	0.31	ng	0.00
8) Nitrobenzene-d5	9.06	82	4154	0.31	ng	0.00
11) 2-Methylnaphthalene-d10	12.29	152	6225	0.23	ng	0.00
14) 2,4,6-Tribromophenol	16.00	330	739	0.36	ng	0.00
15) 2-Fluorobiphenyl	13.15	172	8227	0.21	ng	-0.02
25) Fluoranthene-d10	19.27	212	54069	0.19	ng	0.00
29) Terphenyl-d14	19.86	244	9937	0.22	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
22) Pentachlorophenol	16.91	266	184	0.188	ng	88
23) Phenanthrene	17.27	178	2206	0.040	ng	# 90
26) Fluoranthene	19.30	202	4603	0.053	ng	# 91
28) Pyrene	19.66	202	4304	0.054	ng	# 91
30) Benzo(a)anthracene	21.39	228	2559	0.035	ng	# 87
31) Chrysene	21.45	228	2578	0.032	ng	# 87
32) Bis(2-ethylhexyl)phthalate	21.30	149	18353	0.184	ng	97
33) Indeno(1,2,3-cd)pyrene	26.55	276	1674	0.023	ng	96
35) Benzo(b)fluoranthene	23.15	252	2998	0.038	ng	# 61
37) Benzo(a)pyrene	23.80	252	2164	0.030	ng	# 56
39) Benzo(g,h,i)perylene	27.36	276	1631	0.023	ng	# 78

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

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