

Data Path : Z:\HPCHEM1\BNA E\DATA\BE060816\
 Data File : BE091905.D
 Acq On : 8 Jun 2016 17:34
 Operator : UM/SJ
 Sample : H3359-01RE
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 OM-SPN-001-160531RE

Quant Time: Jun 09 04:58:23 2016
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE060716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 08 16:11:34 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.77	152	30476	5.00	ng	0.00
8) Naphthalene-d8	10.55	136	136901	5.00	ng	0.00
15) Acenaphthene-d10	14.38	164	82921	5.00	ng	0.00
24) Phenanthrene-d10	17.12	188	254735	5.00	ng	0.00
31) Chrysene-d12	21.30	240	204285	5.00	ng	0.00
40) Perylene-d12	23.71	264	282812	5.00	ng	0.00

System Monitoring Compounds						
4) 2-Fluorophenol	5.38	112	25219	3.88	ng	0.00
5) Phenol-d6	6.94	99	35343	4.39	ng	-0.02
9) Nitrobenzene-d5	8.91	82	19542	1.92	ng	-0.02
16) 2,4,6-Tribromophenol	15.87	330	9904	4.19	ng	-0.02
17) 2-Fluorobiphenyl	13.01	172	54679	2.17	ng	0.00
34) Terphenyl-d14	19.74	244	104871	3.16	ng	0.00

Target Compounds						Qvalue
26) Hexachlorobenzene	16.44	284	195	0.02	ng	# 87
27) Pentachlorophenol	16.78	266	221	0.11	ng	# 88
28) Phenanthrene	17.17	178	1607	0.03	ng	95
29) Anthracene	17.26	178	1178m	0.02	ng	
30) Fluoranthene	19.19	202	1631	0.03	ng	97
33) Pyrene	19.53	202	1765	0.03	ng	99
37) Chrysene	21.33	228	6623m	0.03	ng	
38) Bis(2-ethylhexyl)phthalate	21.19	149	10240	0.39	ng	# 100
39) Indeno(1,2,3-cd)pyrene	26.29	276	1840	0.07	ng	# 76
41) Benzo(b)fluoranthene	22.98	252	2060	0.03	ng	# 73
42) Benzo(k)fluoranthene	23.03	252	1995	0.03	ng	# 74
43) Benzo(a)pyrene	23.59	252	1704	0.03	ng	# 51
44) Dibenzo(a,h)anthracene	26.31	278	1436	0.02	ng	# 70
45) Benzo(a,h,i)perylene	27.07	276	1657	0.03	ng	# 81

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

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