

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030316\
 Data File : BM004549.D
 Acq On : 04 Mar 2016 01:37
 Operator : SJ/UM
 Sample : H1616-10
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P009-SS003-01

Manual Integrations
 APPROVED

UMANGI
 3/4/2016 4:21:17 PM

Quant Time: Mar 04 06:41:45 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM030316.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 04 04:25:59 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	38700	20.00	ng/ul	0.00
18) Naphthalene-d8	10.38	136	188169	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.26	164	111008	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.02	188	219289	20.00	ng/ul	0.00
78) Chrysene-d12	21.23	240	153302	20.00	ng/ul	0.00
86) Perylene-d12	23.44	264	150542	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.05	96	2123	2.95	ng/uL	0.00
5) Phenol-d5	6.85	99	60032	19.25	ng/ul	0.06
7) Bis-(2-Chloroethyl)ether-d	6.93	67	28998	18.45	ng/ul	0.00
9) 2-Chlorophenol-d4	7.15	132	55622	20.01	ng/ul	0.02
13) 4-Methylphenol-d8	8.36	113	51515	18.74	ng/ul	0.04
19) Nitrobenzene-d5	8.76	128	29072	20.67	ng/ul	0.00
22) 2-Nitrophenol-d4	9.47	143	33898	21.06	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.07	165	59410	20.40	ng/ul	0.05
29) 4-Chloroaniline-d4	10.53	131	50664	14.14	ng/ul	0.00
44) Dimethylphthalate-d6	13.67	166	191866	20.78	ng/ul	0.00
47) Acenaphthylene-d8	13.95	160	236293	21.16	ng/ul	0.00
52) 4-Nitrophenol-d4	14.66	143	14766m	10.84	ng/ul	0.15
58) Fluorene-d10	15.26	176	165733	20.69	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.42	200	23050	16.19	ng/ul	0.00
71) Anthracene-d10	17.12	188	226651	21.12	ng/ul	0.00
79) Pyrene-d10	19.42	212	197709	23.98	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.30	264	167128	20.87	ng/ul	0.01

Target Compounds

						Qvalue
28) Naphthalene	10.43	128	30625	2.96	ng/ul	99
45) Dimethylphthalate	13.72	163	73773	7.60	ng/ul	99
70) Phenanthrene	17.06	178	31155	2.37	ng/ul	97
77) Fluoranthene	19.09	202	47408	3.35	ng/ul	99
80) Pyrene	19.45	202	50832	4.57	ng/ul	99
83) Benzo(a)anthracene	21.21	228	22157	2.24	ng/ul#	90
84) Bis(2-ethylhexyl)phthalate	21.13	149	31911	5.15	ng/ul#	96
85) Chrysene	21.26	228	23823	2.51	ng/ul#	90
88) Benzo(b)fluoranthene	22.78	252	28962	2.89	ng/ul#	90
91) Benzo(a)pyrene	23.34	252	24792	2.62	ng/ul#	88
92) Indeno(1,2,3-cd)pyrene	25.67	276	18454	1.95	ng/ul	94
94) Benzo(g,h,i)perylene	26.36	276	21871	2.75	ng/ul	95

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

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030316\
Data File : BM004549.D
Acq On : 04 Mar 2016 01:37
Operator : SJ/UM
Sample : H1616-10
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
P009-SS003-01

Manual Integrations
APPROVED

UMANGI
3/4/2016 4:21:17 PM

Quant Time: Mar 04 06:41:45 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM030316.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Mar 04 04:25:59 2016
Response via : Initial Calibration

