

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG031122\
 Data File : BG052657.D
 Acq On : 11 Mar 2022 21:28
 Operator : CG/JU
 Sample : N1886-04DL 5X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 30970DL

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 03/14/2022
 Supervised By : Jagrut Upadhyay 03/14/2022

Quant Time: Mar 12 06:17:25 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG030922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 10 06:37:54 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.171	152	36049	20.000	ng	0.00
21) Naphthalene-d8	11.009	136	138738	20.000	ng	0.00
39) Acenaphthene-d10	14.827	164	91590	20.000	ng	0.00
64) Phenanthrene-d10	17.576	188	189457	20.000	ng	0.00
76) Chrysene-d12	21.888	240	169640	20.000	ng	0.00
86) Perylene-d12	25.325	264	175857	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.704	112	17079	8.185	ng	0.00
7) Phenol-d6	7.325	99	14026	4.644	ng	0.00
23) Nitrobenzene-d5	9.346	82	45231	14.779	ng	0.00
42) 2,4,6-Tribromophenol	16.319	330	26079	20.076	ng	0.00
45) 2-Fluorobiphenyl	13.446	172	99191	14.618	ng	0.00
79) Terphenyl-d14	20.173	244	114863	11.982	ng	0.00
Target Compounds						Qvalue
6) Aniline	7.490	93	10620	3.117	ng	99
10) Phenol	7.355	94	11871	3.974	ng	97
20) 3+4-Methylphenols	8.947	107	109642	40.116	ng	96
27) 2,4-Dimethylphenol	10.174	122	11727m	6.268	ng	
70) Pentachlorophenol	17.241	266	8551	11.859	ng	# 82
84) Bis(2-ethylhexyl)phtha...	21.741	149	17813	3.090	ng	93

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

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