

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021121\
 Data File : BP004714.D
 Acq On : 11 Feb 2021 14:02
 Operator : CG/JU
 Sample : M1354-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 T22-216481

Quant Time: Feb 11 14:39:00 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 05 18:04:48 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.769	152	65466	20.00	ng	0.00
21) Naphthalene-d8	10.563	136	243267	20.00	ng	# 0.00
39) Acenaphthene-d10	14.422	164	155009	20.00	ng	0.00
64) Phenanthrene-d10	17.192	188	343775	20.00	ng	# 0.02
76) Chrysene-d12	21.263	240	319763	20.00	ng	0.00
86) Perylene-d12	23.563	264	317362	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.358	112	60258	14.21	ng	0.00
7) Phenol-d6	6.934	99	71905	12.33	ng	0.00
23) Nitrobenzene-d5	8.922	82	46552	8.34	ng	-0.01
42) 2,4,6-Tribromophenol	15.922	330	25680	12.47	ng	0.00
45) 2-Fluorobiphenyl	13.034	172	98236	7.75	ng	0.00
79) Terphenyl-d14	19.757	244	164098	8.62	ng	0.02
Target Compounds						Qvalue
37) 2-Methylnaphthalene	12.228	142	35351	3.50	ng	# 88
38) 1-Methylnaphthalene	12.452	142	36105	3.78	ng	# 96
46) 1,1'-Biphenyl	13.251	154	53891	4.15	ng	92
55) Dibenzofuran	14.822	168	37443	2.39	ng	# 2
71) Phenanthrene	17.234	178	166328	7.90	ng	# 7

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021121\
Data File : BP004714.D
Acq On : 11 Feb 2021 14:02
Operator : CG/JU
Sample : M1354-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
T22-216481

Quant Time: Feb 11 14:39:00 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Feb 05 18:04:48 2021
Response via : Initial Calibration

