

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110121\
 Data File : BN017223.D
 Acq On : 01 Nov 2021 21:40
 Operator : CG/JU
 Sample : M4360-04
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BP-VPB-RW9-GW-148-150

Quant Time: Nov 02 03:56:47 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN110121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 01 17:00:24 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.561	152	29967	0.400	ng	0.00
7) Naphthalene-d8	10.319	136	138211	0.400	ng	0.00
13) Acenaphthene-d10	14.194	164	75068	0.400	ng	0.00
19) Phenanthrene-d10	16.946	188	148454	0.400	ng	0.00
29) Chrysene-d12	21.155	240	134369	0.400	ng	# 0.00
35) Perylene-d12	23.356	264	135720	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.191	112	10484	0.144	ng	0.02
5) Phenol-d6	6.749	99	7037	0.089	ng	0.00
8) Nitrobenzene-d5	8.697	82	24019	0.283	ng	0.00
11) 2-Methylnaphthalene-d10	11.920	152	68892	0.353	ng	0.00
14) 2,4,6-Tribromophenol	15.691	330	8448	0.324	ng	0.00
15) 2-Fluorobiphenyl	12.811	172	93059	0.330	ng	0.00
27) Fluoranthene-d10	18.985	212	137994	0.365	ng	0.00
31) Terphenyl-d14	19.596	244	117445	0.400	ng	0.00
Target Compounds						
24) Pentachlorophenol	16.605	266	168	0.209	ng	90
34) Bis(2-ethylhexyl)phtha...	21.091	149	188985	0.852	ng	94

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110121\
Data File : BN017223.D
Acq On : 01 Nov 2021 21:40
Operator : CG/JU
Sample : M4360-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BP-VPB-RW9-GW-148-150

Quant Time: Nov 02 03:56:47 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN110121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Nov 01 17:00:24 2021
Response via : Initial Calibration

