

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111320\
 Data File : BP003963.D
 Acq On : 13 Nov 2020 20:06
 Operator : CG/JU
 Sample : L4719-12
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP-3A-S

Manual Integrations
 APPROVED

mohammad
 11/19/2020 10:28:11 AM

Quant Time: Nov 16 08:21:00 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP110320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 12 15:22:47 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.281	152	135040	20.00	ng	0.00
21) Naphthalene-d8	11.110	136	516457	20.00	ng	# 0.00
39) Acenaphthene-d10	14.904	164	308405	20.00	ng	0.00
64) Phenanthrene-d10	17.633	188	638895	20.00	ng	# 0.00
76) Chrysene-d12	21.669	240	575701	20.00	ng	0.00
86) Perylene-d12	24.151	264	573156	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.805	112	914114	112.98	ng	0.00
7) Phenol-d6	7.452	99	1054162	90.76	ng	0.00
23) Nitrobenzene-d5	9.446	82	873428	82.83	ng	0.00
42) 2,4,6-Tribromophenol	16.386	330	490363	127.12	ng	0.00
45) 2-Fluorobiphenyl	13.528	172	1803639	81.76	ng	0.00
79) Terphenyl-d14	20.133	244	2409085	80.85	ng	0.00
Target Compounds						
32) Benzoic acid	10.463	122	6076m	9.35	ng	Qvalue
50) Dimethylphthalate	14.328	163	99421	4.12	ng	98

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111320\
Data File : BP003963.D
Acq On : 13 Nov 2020 20:06
Operator : CG/JU
Sample : L4719-12
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TP-3A-S

Manual Integrations
APPROVED

mohammad
11/19/2020 10:28:11 AM

Quant Time: Nov 16 08:21:00 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP110320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Nov 12 15:22:47 2020
Response via : Initial Calibration

