

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102220\
 Data File : BP003708.D
 Acq On : 23 Oct 2020 05:14
 Operator : CG/JU
 Sample : L4467-02
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 VB-15242

Manual Integrations
 APPROVED

mohammad
 10/23/2020 1:32:53 PM

Quant Time: Oct 23 07:37:46 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP101920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 19 14:54:24 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.528	152	165591	20.00	ng	0.00
21) Naphthalene-d8	11.369	136	617775	20.00	ng	# 0.00
39) Acenaphthene-d10	15.122	164	418810	20.00	ng	0.00
64) Phenanthrene-d10	17.839	188	974710	20.00	ng	0.00
76) Chrysene-d12	21.851	240	1045960	20.00	ng	# 0.00
86) Perylene-d12	24.421	264	963832	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	6.028	112	1185282	119.69	ng	0.00
7) Phenol-d6	7.675	99	1389402	99.31	ng	0.00
23) Nitrobenzene-d5	9.693	82	1240644	84.96	ng	0.00
42) 2,4,6-Tribromophenol	16.598	330	902475	138.12	ng	0.00
45) 2-Fluorobiphenyl	13.751	172	2823647	87.16	ng	0.00
79) Terphenyl-d14	20.310	244	4931118	82.92	ng	0.00
Target Compounds						
10) Phenol	7.699	94	70979	5.33	ng	Qvalue 73
32) Benzoic acid	10.593	122	3488m	5.58	ng	
50) Dimethylphthalate	14.539	163	241163	7.01	ng	99

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

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