

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100720\
 Data File : BP003567.D
 Acq On : 07 Oct 2020 23:23
 Operator : CG/JU
 Sample : L4221-09
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 OK-01-100620

Quant Time: Oct 08 01:55:17 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 05 14:33:14 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.493	152	78569	20.00	ng	0.00
21) Naphthalene-d8	10.269	136	303933	20.00	ng	# 0.00
39) Acenaphthene-d10	14.151	164	176360	20.00	ng	0.00
64) Phenanthrene-d10	16.910	188	347917	20.00	ng	0.00
76) Chrysene-d12	21.016	240	303338	20.00	ng	-0.01
86) Perylene-d12	23.186	264	352656	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.111	112	578320	128.53	ng	0.00
7) Phenol-d6	6.699	99	669621	111.65	ng	0.00
23) Nitrobenzene-d5	8.646	82	559032	108.06	ng	0.00
42) 2,4,6-Tribromophenol	15.657	330	314573	116.95	ng	0.00
45) 2-Fluorobiphenyl	12.763	172	1125176	93.31	ng	-0.01
79) Terphenyl-d14	19.492	244	1405618	96.55	ng	-0.01

Target Compounds	Qvalue

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

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