

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110620\
 Data File : BP003858.D
 Acq On : 07 Nov 2020 04:45
 Operator : CG/JU
 Sample : L4656-09
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SR-L14-L13-SOUTH

Quant Time: Nov 07 05:32:56 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP110320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 06 11:33:14 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.304	152	144768	20.00	ng	0.00
21) Naphthalene-d8	11.134	136	559227	20.00	ng	# 0.00
39) Acenaphthene-d10	14.916	164	345385	20.00	ng	0.00
64) Phenanthrene-d10	17.645	188	717892	20.00	ng	0.00
76) Chrysene-d12	21.662	240	681936	20.00	ng	0.00
86) Perylene-d12	24.145	264	701761	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.828	112	1063611	122.62	ng	0.00
7) Phenol-d6	7.469	99	1234806	99.17	ng	0.00
23) Nitrobenzene-d5	9.463	82	978100	85.66	ng	0.00
42) 2,4,6-Tribromophenol	16.398	330	654683	151.55	ng	0.00
45) 2-Fluorobiphenyl	13.545	172	2176785	88.11	ng	0.00
79) Terphenyl-d14	20.139	244	3330980	94.37	ng	0.00
Target Compounds						
10) Phenol	7.493	94	25036	2.08	ng	# 68
50) Dimethylphthalate	14.339	163	156436	5.79	ng	100

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

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