

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102220\
 Data File : BP003714.D
 Acq On : 23 Oct 2020 08:38
 Operator : CG/JU
 Sample : L4225-14
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SU-02-102120

Quant Time: Oct 23 09:12:29 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	152324	20.00	ng	0.00
21) Naphthalene-d8	11.369	136	572119	20.00	ng	# 0.00
39) Acenaphthene-d10	15.122	164	390008	20.00	ng	0.00
64) Phenanthrene-d10	17.839	188	913111	20.00	ng	0.00
76) Chrysene-d12	21.851	240	1032859	20.00	ng	# 0.00
86) Perylene-d12	24.427	264	991024	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	6.022	112	1109694	121.82	ng	0.00
7) Phenol-d6	7.675	99	1318017	102.41	ng	0.00
23) Nitrobenzene-d5	9.693	82	1136057	84.00	ng	0.00
42) 2,4,6-Tribromophenol	16.598	330	826449	135.83	ng	0.00
45) 2-Fluorobiphenyl	13.751	172	2582319	85.60	ng	0.00
79) Terphenyl-d14	20.310	244	4560653	77.66	ng	0.00
Target Compounds						
10) Phenol	7.699	94	72344	5.90	ng	78
50) Dimethylphthalate	14.539	163	248768	7.76	ng	99

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

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