

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040121\
 Data File : BP005137.D
 Acq On : 01 Apr 2021 15:44
 Operator : CG/JU
 Sample : M1832-10
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 IG005

Quant Time: Apr 01 16:32:40 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP032321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 01 12:23:57 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.705	152	25558	20.00	ng	# 0.00
21) Naphthalene-d8	10.493	136	95947	20.00	ng	# 0.00
39) Acenaphthene-d10	14.363	164	62789	20.00	ng	0.00
64) Phenanthrene-d10	17.122	188	133291	20.00	ng	# 0.00
76) Chrysene-d12	21.216	240	143366	20.00	ng	-0.01
86) Perylene-d12	23.504	264	149941	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.305	112	228689	137.70	ng	0.00
7) Phenol-d6	6.887	99	262252	110.24	ng	0.00
23) Nitrobenzene-d5	8.863	82	218768	97.83	ng	0.00
42) 2,4,6-Tribromophenol	15.857	330	119720	146.34	ng	0.00
45) 2-Fluorobiphenyl	12.975	172	428496	92.32	ng	0.00
79) Terphenyl-d14	19.698	244	729683	92.65	ng	0.00
Target Compounds						
10) Phenol	6.911	94	4934	2.02	ng	# 69
50) Dimethylphthalate	13.822	163	9763	2.02	ng	# 96

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

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