

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP073021\
 Data File : BP006446.D
 Acq On : 30 Jul 2021 20:08
 Operator : CG/JU
 Sample : PB137988BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB137988BL

Quant Time: Jul 31 00:42:37 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270-SIM-BP073021.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 30 18:59:15 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.785	152	4439	0.400	ng	0.00
7) Naphthalene-d8	10.546	136	16776	0.400	ng	0.00
13) Acenaphthene-d10	14.390	164	8015	0.400	ng	0.00
19) Phenanthrene-d10	17.135	188	16038	0.400	ng	0.00
29) Chrysene-d12	21.253	240	13687	0.400	ng	0.00
35) Perylene-d12	23.583	264	12683	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.357	112	4277	0.333	ng	0.00
5) Phenol-d6	6.940	99	6549	0.348	ng	0.00
8) Nitrobenzene-d5	8.912	82	6400	0.354	ng	0.00
11) 2-Methylnaphthalene-d10	12.137	152	8672	0.370	ng	0.00
14) 2,4,6-Tribromophenol	15.880	330	1127	0.305	ng	0.00
15) 2-Fluorobiphenyl	13.024	172	12387	0.369	ng	0.00
27) Fluoranthene-d10	19.140	212	16408	0.370	ng	0.00
31) Terphenyl-d14	19.729	244	11188	0.380	ng	0.00

Target Compounds Qvalue

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

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