

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082321\
 Data File : BP006830.D
 Acq On : 23 Aug 2021 14:38
 Operator : CG/JU
 Sample : M3463-43
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 2021-08-17-FD-05

Quant Time: Aug 23 15:34:18 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 23 11:39:38 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.693	152	143869	20.000	ng	0.00
21) Naphthalene-d8	10.481	136	597044	20.000	ng	# 0.00
39) Acenaphthene-d10	14.339	164	345606	20.000	ng	0.00
64) Phenanthrene-d10	17.104	188	717298	20.000	ng	# 0.00
76) Chrysene-d12	21.210	240	703182	20.000	ng	# 0.00
86) Perylene-d12	23.527	264	755211	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.305	112	1386531	148.026	ng	0.00
7) Phenol-d6	6.887	99	1732059	130.174	ng	0.00
23) Nitrobenzene-d5	8.857	82	1165502	90.105	ng	0.00
42) 2,4,6-Tribromophenol	15.845	330	622906	172.456	ng	0.00
45) 2-Fluorobiphenyl	12.963	172	2123990	91.947	ng	0.00
79) Terphenyl-d14	19.686	244	3100467	88.351	ng	0.00

Target Compounds	Qvalue

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

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