

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF100821\
 Data File : BF125785.D
 Acq On : 8 Oct 2021 16:28
 Operator : JU/CG
 Sample : M4087-10
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-20211005-ROCK

Quant Time: Oct 08 17:27:10 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF100721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 07 17:16:23 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.845	152	214685	20.00	ng	0.00
21) Naphthalene-d8	8.122	136	860988	20.00	ng	# 0.00
39) Acenaphthene-d10	9.875	164	446977	20.00	ng	0.00
64) Phenanthrene-d10	11.357	188	729277	20.00	ng	# 0.00
76) Chrysene-d12	13.992	240	419280	20.00	ng	0.00
86) Perylene-d12	15.445	264	449532	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.469	112	1372200	104.42	ng	0.01
7) Phenol-d6	6.469	99	1506488	89.05	ng	0.00
23) Nitrobenzene-d5	7.404	82	1087997	78.55	ng	0.00
42) 2,4,6-Tribromophenol	10.663	330	492408	111.58	ng	0.00
45) 2-Fluorobiphenyl	9.198	172	2098647	80.98	ng	0.00
79) Terphenyl-d14	12.945	244	1916997	69.64	ng	0.00

Target Compounds	Qvalue

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

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