

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080122\
 Data File : BF129474.D
 Acq On : 01 Aug 2022 15:30
 Operator : CG\JU
 Sample : N3977-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MS-TOTE-0729

Quant Time: Aug 02 01:16:26 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 02 01:11:25 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.875	152	172906	20.000	ng	0.00
21) Naphthalene-d8	8.157	136	672176	20.000	ng	0.00
39) Acenaphthene-d10	9.910	164	334639	20.000	ng	0.00
64) Phenanthrene-d10	11.404	188	474685	20.000	ng	# 0.00
76) Chrysene-d12	14.045	240	351639	20.000	ng	0.00
86) Perylene-d12	15.533	264	396981	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	517964	48.799	ng	0.00
7) Phenol-d6	6.522	99	386625	28.979	ng	0.00
23) Nitrobenzene-d5	7.440	82	1051496	85.394	ng	0.00
42) 2,4,6-Tribromophenol	10.710	330	375042	116.570	ng	0.00
45) 2-Fluorobiphenyl	9.234	172	1815514	83.926	ng	0.00
79) Terphenyl-d14	12.992	244	1381014	74.093	ng	0.00

Target Compounds Qvalue

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

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