

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080222\
 Data File : BM036071.D
 Acq On : 03 Aug 2022 02:59
 Operator : CG/JU
 Sample : N3894-06
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 RVE-25

Quant Time: Aug 03 03:33:44 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 01 17:50:14 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.869	152	381603	20.000	ng	0.00
21) Naphthalene-d8	10.675	136	1642472	20.000	ng	0.00
39) Acenaphthene-d10	14.504	164	959733	20.000	ng	0.00
64) Phenanthrene-d10	17.245	188	1943102	20.000	ng	0.00
76) Chrysene-d12	21.421	240	1535805	20.000	ng	0.00
86) Perylene-d12	23.791	264	1295454	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.434	112	3491917	148.356	ng	0.00
7) Phenol-d6	7.045	99	4445016	148.416	ng	0.00
23) Nitrobenzene-d5	9.039	82	2760546	94.082	ng	0.00
42) 2,4,6-Tribromophenol	15.998	330	1798270	159.719	ng	0.00
45) 2-Fluorobiphenyl	13.139	172	6182510	95.113	ng	0.00
79) Terphenyl-d14	19.874	244	9021903	115.869	ng	0.00
Target Compounds						
78) Pyrene	19.662	202	271933	2.855	ng	98
88) Benzo(b)fluoranthene	23.068	252	173594	2.287	ng	# 75

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

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