

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080522\
 Data File : BM036133.D
 Acq On : 05 Aug 2022 23:40
 Operator : CG/JU
 Sample : N3961-07
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 RVE-174

Quant Time: Aug 06 01:14:22 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 05 03:36:25 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.857	152	283076	20.00	ng	0.00
21) Naphthalene-d8	10.663	136	1171772	20.00	ng	# 0.00
39) Acenaphthene-d10	14.498	164	681275	20.00	ng	0.00
64) Phenanthrene-d10	17.239	188	1351068	20.00	ng	0.00
76) Chrysene-d12	21.415	240	1227384	20.00	ng	0.00
86) Perylene-d12	23.774	264	1272927	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.446	112	1874989	107.39	ng	0.02
7) Phenol-d6	7.069	99	2318239	104.35	ng	0.04
23) Nitrobenzene-d5	9.028	82	1385857	66.20	ng	0.00
42) 2,4,6-Tribromophenol	15.986	330	890577	111.43	ng	0.00
45) 2-Fluorobiphenyl	13.122	172	3160540	68.50	ng	0.00
79) Terphenyl-d14	19.862	244	4497079	72.27	ng	0.00
Target Compounds						Qvalue
75) Fluoranthene	19.292	202	254898	3.31	ng	98
78) Pyrene	19.656	202	294653	3.87	ng	100
88) Benzo(b)fluoranthene	23.056	252	226175	3.03	ng	# 87
90) Benzo(a)pyrene	23.668	252	164672	2.63	ng	# 91

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

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